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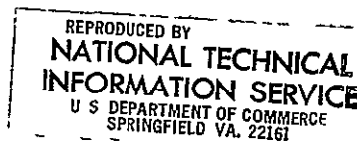
STICAP

A LINEAR CIRCUIT ANALYSIS PROGRAM WITH STIFF SYSTEMS CAPABILITY
Volume III - Systems Programmer's Manual

by Charlie H. Cooke and M. Niel Ransom

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OLD DOMINION UNIVERSITY
School of Engineering
Norfolk, VA 23508

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ABSTRACT

STICAP (Stiff Circuit Analysis Program) is a FORTRAN IV, Version 2.3, computer program written for the CDC-6400-6600 computer series and SCOPE 3.0 operating system. It provides the circuit analyst a tool for automatically computing the transient responses and frequency responses of large linear time invariant networks, both stiff and non-stiff. The circuit description and user's program input language is engineer-oriented, making simple the task of using the program.

Three volumes of documentation are available for the STICAP program; a theory manual, a user's manual, and a system's programmers manual. Volume I describes the engineering theories underlying STICAP and gives further references to the literature. Volume II, the user's manual, explains user interaction with the program and gives results of typical circuit design applications. Volume III depicts the program structure from a system's programmers viewpoint and contains flow charts and other software documentation.

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CHAPTER I

INTRODUCTION

This document contains material describing the internal program structure of the STICAP program. An overview of the composite system structure is given. The overlay structure is described, and the routines composing each overlay are included, for each routine generated under contract NAS1-9434-25. A description of the CORNAP and Gear routines may be found in the references [1] and [2] below, which contain as much documentation as is available for these programs.

The program is written using CDC FORTRAN 4, Version 2.3. It is machine compatible with the CDC 6400-6600 computer series, and runs under the SCOPE 3.0 operating system. It is segmented in overlays of 70K or less using the SCOPE OVERLAY capability. All I/O is accomplished using standard I/O files. I/O files are equivalenced so that File 5 is used for input and File 6 for output. No other files are used by this program.

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1. Pottle, C., "Program Cornap - FORTRAN Computer Routine," System Theory Group, School of Electrical Engineering, Cornell University, Ithaca, N. Y., 1968.
 2. Gear, C. W., Algorithm 407 - DIFSUB for Solution of Ordinary Differential Equations [D2], Communications of the ACM, March, 1971, Vol. 14, No. 3, 185-190.

CHAPTER II

OVERLAY STRUCTURE

The first component of the program overlay structure (see Figure 1) consists of a root segment (a main overlay) containing a supervisor routine STICAP, a commonly used routine USEFCN, and blocks of labeled and unlabeled common core storage, also used by many routines. The supervisor routine obtains control at initial execution and coordinates the calling into core of the successive overlays. This is accomplished by means of subsequent primary overlays, of which each successive one replaces the preceding, with the number called dependent upon the mode of operation selected by the user.

The first primary overlay consists of a block of routines which read the input circuit description cards, scaling card, and mode card. The circuit description cards, whose format has been altered from that of CORNAP, provide the user an easy means of describing the circuit. The scaling card, optionally present, determines the way the circuit equations are to be scaled for numerical computation of a solution. The mode card determines the program mode, or method by which the solution is to be obtained.

The second primary overlay consists of the original program CORNAP, minus the data input processing routines, whose function has been accomplished by the routines mentioned above. The routines of this overlay process the circuit description data previously read, to obtain the mathematical equations governing the circuit. If the solution is to be obtained by the CORNAP mode, the original CORNAP routines continue processing the circuit equations.

In the event that either the Gear or Matrix mode is selected, the third primary overlay is called into core. Data compiled by

Root Segment		STICAP					
		USEFCN					
		LABELED COMMON					
		BLANK COMMON					
Primary Overlays							
PEED	C	COMPAK	INTGR	MATRIX			Secd. Jary Overlays
INPUT	O	APACK*	DIFSUB**	SETSVE	INTCL	OUTPNT	
MODE	R	CROLCR	DATPNT	VANINV	SVECOS	CLSOL	
APACK	N	SRCORD	DIFFUN	ELGINV	SVESIN	TABPNT	
ATODEC	A	RUNCTL	PEDERV	HESFT	SVESTP	ATAN3	
NAME	P	ATODEC*	DECOIP**	ORTRAN	INIVAL		
NODE		DECEQV*	SOLVE**		MPROD		
DECEQV		CROLAX	4th Primary overlay optionally chosen		RAT		
OUTPUT		SRCINT			SCFYAL		
EFSCAL		RUNCRL			TRANS		
		CROLCR			STEADY		
1	2	3			STORE		

Figure 1 - Overlay Structure

*Appears in a previous overlay also.

**Gear's ALGORITHM 407 Routines.

CORNAP is compacted into a more economical form by the routine COMPAK. The control cards and data input cards for the mode selected are processed, and the final primary overlay is called.

The final overlay routines are determined by the mode selection. In one case the routines of Gear are used for numerical integration; in the other the matrix technique is used to obtain the solution.

CHAPTER III

SYSTEM OVERVIEW

An overview of the STICAP system indicacing the overall logical flow of the program appears in the following pages. The figures indicate the routines within each overlay; the functions performed; and the calling sequences, or flow of control.

5

- 6 -

Overlay (1,0)

START

REED

Read and interpret header cards

INPUT

Read and interpret elements card group

OUTPUT

Read and interpret output card group

MODE

Read and interpret mode card

SCALE

Read and interpret scaling card group

Overlay (2,0)

CORNAP I

Generate state equations

CORNAP MODE ?

no

yes

CORNAP II

Solve using CORNAP

END

Overlay (3,0)

COMPAK

Rearrange Blank Common

RUNCTL

Read and interpret RUN CONTROL card group for GEAR

RUNCTL

Read and interpret RUN CONTROL card group for MATRIX mode

which mode

GEAR

MATRIX

CROLGR

Interpret header cards for GEAR mode

CROLMX

Interpret header cards for MATRIX mode

α

ICONS

Read and interpret initial conditions card group

SRCORD

Read and interpret source order card group

SRCINT

Read and interpret source DEFINITION card group

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CHAPTER IV
FUNCTIONS OF THE INDIVIDUAL ROUTINES

- 8 -

In this section a brief description of the function performed by each contractor supplied routine is given, and the routines are categorized by the overlay in which they operate. We remark that occasionally a routine may be used in several overlays; hence it may appear more than once. For a more complete description of a routine, the next section may be consulted.

The name and function of the various routines will now be given.

I. Root Segment Routines

1. USEFCN - Supplies the user a means of specifying the independent sources of the circuit.
2. STICAP - Supervisor routine which coordinates the calling into core of the primary overlays.

II. First Primary Overlay

1. REED - Processes title card and scaling card; coordinates routines which read CORNAP circuit description.
2. INPUT - Reads circuit description cards in a free format.
3. MODE - Reads Mode select card and processes it.
4. APACK - Packs into one word the first character of each word from an array of words.
5. ATODEC - Interprets free format floating point numbers.
6. NAME - Scan routine for breaking out a name from a data field of free format data.

*Some routines not contractor supplied are also indicated. Such routines are Langley routines or else Gear supplementary routines.

7. NODE - Scan routine for breaking out an integer number from a data field of free form data.
8. DECEQV - Changes alphanumeric BCD characters into internal integer format on the CDC - 6600 computer.
9. OUTPUT - Reads and processes cards indicating the circuit outputs specified by the user.
10. ZFSCAL - Reads scaling cards and stores scaling parameters.

III. Second Primary Overlay

CORNAP - The body of routines composing the original program CORNAP, minus routines which read into core the circuit description cards. This overlay also contains the logical decisions required for carrying out the various print options, such as state equations, transfer functions, etc.

IV. Third Primary Overlay

1. APACK - Previously described; first primary overlay.
2. CROLGR - Coordinates the subroutines which read the source order; initial conditions, run controls card group, and end card for the Gear mode.
3. ICONSV - Reads and processes the initial conditions card groups for either the Gear or Matrix modes.
4. SRCORD - Reads and processes the source order card groups for the Gear modes.
5. RUNCTL - Reads and processes run control cards for Gear mode.
6. ATODEC - Previously described; first primary overlay.

7. DECEQV - Previously described: first primary overlay.
8. CROLLK - Coordinates the subroutines which read the source definition, initial conditions, run controls card groups, and the END card for the Matrix mode.
9. SRCINT - Interprets the source definition card group for the Matrix mode.
10. RUHCRL - Reads and processes run controls card groups for the Matrix mode.
11. COMPAK - Rearranges data tables output by CORNAP circuit translation routines into a more economical form.
12. CROLLM - Controlling routine for third overlay.

V. Gear Primary Overlay

1. INTGR - Coordinates Gear's routine DIFSUB for numerical integration and determines when the stop condition is reached.
2. DIFSUB - Integrates the circuit equations in the Gear mode.
3. DATPNT - Prints output data.
4. DIFFUN - Evaluates the right members of the state variable differential equations of the circuit, given a specified time point.
5. PEDERV - Evaluates the Jacobian Matrix of the system equation for the circuit.
6. DECOMP - LU decomposition routine used in tandem with SOLVE to solve a linear system.
7. SOLVE - Second stage of a LU decomposition linear equation solver pair of routines, used by DIFSUB.

VI. Matrix Primary Overlay

1. MATRIX - Calls matrix mode secondary overlays.
2. SETSVE - A controlling routine for matrix mode, which also obtains a sparse matrix code.
3. VANINV - Inverts a Vandermonde matrix.
4. EIGNV - Obtains system eigenvalues using QR transformation.
5. HESEN - Reduces a matrix to upper Hessenburg form.
6. QRTRAN - Performs a single double-step of the QR transformation, avoiding complex algebra.
7. INTCL - Controls the closed form integration performed by the matrix mode.
8. SVECOS - Processes integration of COSINE inputs.
9. SVESIN - Processes integration of SINE inputs.
10. SVESTP - Processes integration of STEP inputs.
11. INIVAL - Integrates for impulse driving functions or other driving functions with a specified initial value.
12. MPROD - Forms a matrix product.
13. EAT - Calculates the matrix exponential $\exp(AT)$ and multiplies into a vector.
14. SCEVAL - Evaluates the $D\bar{u}$ vector, \bar{u} the vector of independent source inputs, $\bar{y} = C\bar{x} + D\bar{u}$ the requested output vector, at a specified time.
15. TRANS - Evaluates transient closed form solution at a specified time point.
16. STEADY - Evaluates steady state closed form solution at a specified time point.
17. STORE - Stores symbolic steady state closed form solution.

- 18. OUTPNT - Controls printing of matrix mode outputs.
- 19. CLSOL - Prints symbolic closed form solutions.
- 20. TABPNT - Prints tabular form of matrix mode solutions.
- 21. ATAN3 - Combines a sine and cosine function to obtain a single sine function.

Subroutine: USEFCN

This routine allows the user the capability of defining the form of the input vector for state-input and state output equations of the network, for the Gear node:

$$\frac{d\bar{x}}{dt} = A\bar{x} + B\bar{u}$$

where

\bar{x} is the state vector of the circuit,

\bar{u} is the vector of independent sources,

\bar{y} is the vector of outputs requested by the user.

The user must either furnish FORTRAN equations defining the independent sources, or else must supply a routine for computing their values, at a given time. The manner in which this may be accomplished is described in the user's manual.

The net effect of an entry to USEFCN must be the following: Given an initial time T and a vector array U, compute values of the independent sources at time T and store as the components of the vector U. The order in which these components are stored must correspond to the order specified by the user when assembling the source order cards group (see user's manual).

Inputs: a value of T and a vector array U

Outputs. The values of the independent sources at time T, stored in the array U, in the order specified by the user's source order cards.

Program IZED

This program reads the title card and calls INPUT, OUTPUT, and ZFSCAL to read the elements, output, and scaling card groups. The card groups are signaled by the following header cards:

*ELEMENTS

*OUTPUTS

*SCALING

where the asterisks appear in column one. The element and outputs card groups must always appear. The scaling card group need be present only if the circuit is to be scaled. The outputs card group must not be given before the elements card group.

This program then calls MODE to interpret the mode card. The mode card must appear after the above card groups.

The outputs from this routine are described in the comment cards preceding the subroutines INPUT, OUTPUT, and ZFSCAL.

SUBROUTINE INPUT

This subroutine reads the elements card group. All cards read until the next card with an asterisk in column 1 are considered element cards.

The cards are of the form

NAME N1 N2 VALUE

for passive elements or independent sources and

NAME N1 N2 VALUE*NAME2

for dependent sources. NAME, N1, N2, and VALUE must be separated by at least one blank. The asterisk between VALUE and NAME2 need not be present. NAME is the user name for the circuit element. This name

must be no more than 4 characters long. The first letter of the element name tells the type of element as follows.

First Letter	Element Type
V	Voltage Source
I	Current Source
R	Resistor
L	Inductor
C	Capacitor
K	Coefficient of Coupling
M	Mutual Inductance or Capacitance

N1 and N2 for non-mutual elements are the number of the nodes between which the element is connected. The circuit nodes should be numbered compactly from zero, although the failure to do so is non-fatal.

N1 is given a positive reference with respect to N2. For mutual elements N1 and N2 are the name of the two elements involved. For dependent sources NAME2 is the name of the controlling element with a V or an I added to the beginning of the name depending on whether the source is voltage or current controlled.

A feature not normally employed by the average user is the following:

A capacitor may be forced into the -proper tree- by placing the word tree after the description of the element.

NAME N1 N2 VALUE TREE

An inductor may be forced into the cotree by similarly placing the word cotree after the element description

NAME N1 N2 VALUE COTREE

OUTPUT PARAMETERS

LIST A 3 dimensional matrix whose I,J,K element tells information J of the Ith element of type K.

K	Type
1	Dependent Voltage Source
2	Independent Voltage Source
3	Zero Valued Resistor or Inductor
4	Tree Forced Capacitor
5	Capacitor
6	Resistor
7	Inductor
8	Cotree Forced Inductor
9	Zero Valued Capacitor
10	Independent Current Source
11	Dependent Current Source
12	Mutual Element
J	Information
1	User Name for Element
2	Element Value
3	Zero (later used to define outputs)
4-5	For non-mutual elements the nodes between which the element is placed
4-5	For mutual elements the names of the two elements involved

NF A vector whose Ith element is the number of elements of type I (defined as in LIST).

CONBCH An array having two columns whose Ith row is the name of the element which controls the Ith dependent

source. Column 1 is for dependent voltage sources and column 2 is for dependent current sources.

ICON An array constructed like CONECH, but whose elements tell the type of control of the dependent sources, 1 for voltage controlled and 2 for current controlled.

NODEG A vector whose Ith element tells the number of elements and sources connected to the Ith node.

NEX The number of nodes to which no elements are connected.

NODES The largest numbered node given by the user.

Subroutine: MODE (IMODE, IPRINT)

Description: This subroutine reads the mode card to determine what program mode is selected and what print options are chosen.

Program execution will be terminated if an error occurs in describing the program mode. An error in describing a print option is non-fatal, but the corresponding print request will be ignored.

Input Parameters: Mode Card - This card will be in the following form

*MODE NAME, OPTION 1, OPTION 2, OPTION 3

The card is in free form, except for an asterisk in column one. The absence of the asterisk is non-fatal. MODE NAME is the name of the solution program and is one of the following

GEAR

CORVAP

MATRIX.

These names may be abbreviated or misspelled as long as the first letter of the first word is correct.

The options following the name of the solution program are print requests. They are separated by commas and are of the following form:

SOLUTION EQUATIONS

STATE EQUATIONS

TRANSFER FUNCTIONS

These names may be abbreviated or misspelled, as long as the first two letters are correct. The order in which the print options appear is insignificant, nor do all or any need be present.

Output Parameters: IMODE - An integer telling what solution program to use

IMODE = 1 CORNAP

IMODE = 2 GEAR

IMODE = 3 MATRIX

IPRINT - A vector telling what print options were requested.

IPRINT(I) is 1 if option I was requested and 0 if it was not. The elements of IPRINT relate the following:

IPRINT(1) State equations

IPRINT(2) Transfer functions

IPRINT(3) Solution equations

Additional Subroutines Required

APACK

Examples of Usage

1. Card read:

*GEAR MODE

The subroutine would return

```
IMODE = 2  
IPRINT(1) = 0  
IPRINT(2) = 0  
IPRINT(3) = 0
```

2. Card read

```
*MATRIX, SOLUTION EQUATIONS, TRANSFER FUNCTIONS
```

The subroutine would return

```
IMODE = 3  
IPRINT(1) = 0  
IPRINT(2) = 1  
IPRINT(3) = 1
```

3. Card read

```
CORNAP, TRANSFER FUNCTIONS
```

Here the asterisk in column 1 was omitted. The subroutine would return

```
IMODE = 1  
IPRINT(1) = 0  
IPRINT(2) = 1  
IPRINT(3) = 0
```

4. Card read

```
'C,ST EQU,TR
```

Here the names of the solution program and the print requests were abbreviated. The subroutine would return

```
IMODE = 1  
IPRINT(1) = 1  
IPRINT(2) = 1  
IPRINT(3) = 0
```

Subroutine: APACK(NLET,ARRAY,NWORD)

Purpose: To pack the first character of an array of words into one word.

Input Parameters: ARRAY - An array of words containing alpha-numeric BCD characters.

 NLET - The number of words in ARRAY

Output Parameters: NWORD - A word containing the first character of the first word in ARRAY as its first character, the first character of the second word in ARRAY as its second character, and so on until ARRAY is exhausted. The remainder of NWORD is filled with blanks.

Machine Dependency: This subroutine assumes 60 bit words with 6 bit alpha-numeric characters stored from the left.

Example of Usage: ARRAY(1) = 3HABC
 ARRAY(2) = 1HD
 ARRAY(3) = 4HEFGH
 CALL APACK(3,ARRAY,NWORD)

Upon return from the subroutine call, NWORD would contain 3HADE.

Subroutines Required:

 RSHFT

 system subroutines; not STICAP routines.

 RSHFTA

Subroutine: ATODEC (NUFFER, LSRT, LSTP, DEC, NERR)

Description: This subroutine will interpret a number which was read

from a card and stored in an array in alpha-numeric BCD characters.

The number may be in integer, decimal or exponential form. The numbers and its exponent are presumed positive if no sign is given. Numbers in exponential form need not contain a decimal point.

All blanks within the input field are ignored.

The subroutine starts at some given point in the input field and interprets until the first character not part of the number is found. This does not necessarily mean a non-numeric character. The location of this character and the number found is outputted by the subroutine. If no number is found before a non-numeric character, such as a digit or a + or - sign, the number being interpreted is assumed to be a coefficient of 1.0. If no number is found before the end of the input field, a 0.0 is given by the subroutine as the number.

Limitations: The magnitude of the number found must be between 10^{-290} and 10^{+290} and contain no greater than 15 digits.

Input Parameters:

NUFFER - A vector containing the card read in 80A1 format.

LSRT - A starting location in NUFFER where interpreting is to begin.

Output Parameters:

DEC - The number found.

LSTP - The location in NUFFER of the next non-blank character following the number found. If the rest of the card was blank, LSTP = 81

NERR - Error Code

NERR = 0 No error present

NERR = 1 Number is more than 15 digits long
NERR = 2 Number is not between 10^{-290} and 10^{+290} in
 magnitude.

Machine Dependency: There is no machine dependency within the
subroutine itself, however a function which it calls, DECEQV, is
machine dependent.

Examples of Usage:

```
READ 10, (NUFFER(I), I=1, 80)
```

```
10 FORMAT (80 A1)
```

(1) Card contains bbb2.147bbb...

```
CALL ATODEC(NUFFER,1,LSTP, DEC,NERR)
```

After the call

NUFFER remains unchanged

LSTP contains 81

DEC contains 2.147

NERR contains 0

(2) Card contains bbA=2SIN(4T)bb...

```
CALL ATODEC(NUFFER,5,LSTP,DEC,NERR)
```

After the call

NUFFER remains unchanged

LSTP contains 6

DEC contains 2.0

NERR contains 0

(3) Card contains bb216E-3EXP(-X)bb...

```
CALL ATODEC(NUFFER,1,LSTP,DEC,NERR)
```

After the call

NUFFER remains unchanged
 LSTP contains 9
 DEC contains .216
 NERR contains 0

- (4) Card contains -SIN3Zbbb
 CALL ATODEC(NUFFER,1,LSTP,DEC,NERR)

After the call

NUFFER remains unchanged
 LSTP contains 2
 DEC contains -1.0
 NERR contains 0

- (5) Card contains bb3.1415926535897932bb...
 CALL ATODEC(NUFFER,1,LSTP,DEC,NERR)

After the call

NUFFER remains unchanged
 LSTP undefined
 DEC undefined
 NERR contains 1

- (6) Card contains bb-2.14E-5bb21SIN(X)bbb...
 CALL ATODEC(NUFFER,1, LSTP, DEC,NERR)

After the call

NUFFER remains unchanged
 LSTP undefined
 DEC contains -2.14
 NERR contains 2

Subroutines required:

NCODE, DECEOV

Subroutine:

NAME (NUFFER, LSRT, LSTR, NERR)

This subroutine scans a card image to find an element name.

This name must be less than 4 characters long.

Input Parameters:

NUFFER The card image in 80A1 format.

LSRT Starting location where scanning is to begin

Output Parameters:

NWORD The element name found.

LSTP Location of blank after the element name.

NERR Error code.

NERR = 0 No error present.

NERR = 1 Name over 4 characters long.

NERR = 2 No name present (rest of card blank).

Additional Subroutines Required:

APACK

Subroutine:

NODE (NUFFER, LSRT, LSTP, NUM, NERR)

This subroutine scans a card image from a given starting location to find the node number. The subroutine requires this node number to be followed by a blank. The largest node number must be less than 64.

Input Parameters:

NUFFER An array containing the card image in 80A1 format.
LSRT The starting location where scanning is to begin.

Output Parameters:

NUM The node number found.

LSTP The location of the first blank following the
 node number.

NERR Error code

 NERR = 0 No error present

 NERR = 1 A blank does not follow the number.

 NERR = 2 Node number greater than 64.

 NERR = 3 Card after LSRT blank.

Function: DECEQV (NCOEF, LCNT)

Function Purpose: Convert an integer whose digits are BCD characters to an equivalent floating point number.

Limitations: The integer being converted must be less than 15 digits long.

Input Parameters:

NCOEF - An integer array whose Ith element is the Ith digit of the integer being converted, read from the left in A format.

LCNT - The number of digits in the number being converted.

Output Parameters:

DECEQV - The equivalent floating point number.

Machine Dependency: This function assumes 60 bit words with alphanumeric characters in CDC 6000 series BCD code stored on the left.

Example of Usage:

NCOEF (1) = 1H2

NCOEF (2) = 1H8

NCOEF (3) = 1H4

NCOEF (4) = 1H5

DEC = DECEQV (NCOEF, 4)

Upon return from the subroutine call DEC would contain 2845.0.

Subroutines Required:

RSHT (a system subroutine; not a STICAP routine)

Subroutine: OUTPUT

This subroutine reads the output card group to determine the element currents and voltages selected as circuit outputs. This

subroutine will terminate circuit processing if no outputs are selected or if all outputs contain errors.

All cards read until a card with an asterisk in column 1 is found are considered output cards. The cards are of the form

VNAME

or

INAME

depending on whether the voltage across or the current through NAME is desired. Here NAME is the name of a circuit element given in the elements card group.

Input Parameters:

NF - A vector of length 12 whose elements tell the number of each of the 12 types of elements contained in the circuit.

LIST - A matrix whose (I,1,J) element tells the name of the Ith element of type J.

Output Parameters:

LIST - A matrix whose (I,4,J) element gives the output information for the Ith element of type J.

LIST (I,3,J)=0 No output taken from this element.

LIST (I,3,J)=1 Voltage across this element is an output.

LIST (I,3,J)=2 Current through this element is an output.

Subroutine: ZFSCAL

This subroutine reads the scaling card group. The circuit will be scaled around a frequency or impedance value if one of the

following cards are present

FREQUENCY = A

IMPEDANCE = A

where A is a floating point number in free format. Scaling requests containing errors are ignored.

Input Parameters: None

Output Parameters:

ZSCALE - The impedance around which the circuit is to be scaled.

FSCALE - The frequency around which the circuit is to be scaled.

NUFFER - A vector containing the next header card.

Subroutine: COMPAK (EPS,KA,KSC,NDERIV,NROWA,NLC,INODE,IPRINT,FSCALE,ZSCALE,NAIILC)

Description: This subroutine takes information assembled by CORNAP subroutines and puts it in the form used by the GEAR and MATRIX mode subroutines.

This subroutine will terminate execution whenever derivatives of the inputs are required by the output equations or whenever derivatives of the inputs occur higher than the first degree.

Whenever the only derivatives of the inputs appear in the state variable equations and then only the first degree, the state variable and output equations will be of the form

$$\dot{\underline{X}} = \underline{A}\underline{X} + \underline{B}\underline{u} + \underline{B}_1\dot{\underline{u}}$$

$$\underline{Y} = \underline{C}\underline{X} + \underline{D}\underline{u}$$

The following substitutions are then made

$$\underline{A} \rightarrow \underline{A} + \underline{B}_1 + \underline{B} \rightarrow \underline{B}$$

$$\underline{C} \rightarrow \underline{C} + \underline{B}_1 + \underline{D} \rightarrow \underline{D}$$

The new circuit equations now take the form

$$\dot{\underline{q}} = \underline{Aq} + \underline{Bu}$$
$$\underline{Y} = \underline{Cq} + \underline{Du}.$$

The names of the circuit outputs are augmented by adding an I or V in front of the old name depending on whether the current through this element or the voltage across this element is the output.

Inputs and Outputs: All inputs are brought in by the two arrays BLOCK and NAMES stored in common and by the argument list of the subroutine.

The following variables are inputs to the subroutine but are not outputted.

KA,KSC,NROWA -- Constants used to locate the A, B, B1, C, D,
and D1 matrices in BLOCK

D1 - Coefficient matrix of source derivatives in the
output equations

The storage location of the remaining input and output variables are given in Figure 1.

Machine Dependency: The process of adding an I or V in front of the output names is machine dependent.

The addition of the I or V is a means of more readily identifying current and voltage variables for printout and/or scaling purposes.

FIGURE 1

Variable	Length	Starting Location When Loaded	Output Storing Location in MEMORY	Meaning
A	900	BLOCK (KAMNO, $\frac{NRQTA}{2}$, KSC)	1	System Matrix
B	300	BLOCK (KATTO, $\frac{NRQTA}{2}$, KSC, NSTV);	101	Source Coef. Matrix in State eqns
C	300	BLOCK (KA, $\frac{NRQTA}{2}$ + KSC)	1201	State Variable Coef Matrix in Output eqns
D	100	BLOCK (KA, $\frac{NRQTA}{2}$, KSC, NSTV)	1501	Source Coef Matrix in Output eqns
PI	300	BLOCK (KAMNO, $\frac{NRQTA}{2}$, KSC, NSTV, KSC, NIN)	1601	Coef Matrix of Source Deriv. in State eqns
NSTV	1	NAMES (17)	1901	Number of State Variables
NIN	1	NAMES (17)	1902	Number of inputs
NO	1	NAMES (17)	1903	Number of outputs
NLC	1	Subroutine argument	1904	Number of coils and capacitors
NDERIV	1	Subroutine argument	1905	Number of source derivatives
IMODE	1	Subroutine argument	1906	Solution Method, 2-GEAR, 3-MATRIX
IPRINT	3	Subroutine argument	1907	Extra Printouts desired
NAMIN	10	NAMES (49)	1910	Names of inputs
NAMSV	30	NAMES (33)	1920	Names of state variables
NAMOUT	10	NAMES (65)	1950	Names of outputs
NAMLC	30	Subroutine argument	1960	Names of coils and capacitors
FSCALE	1	Subroutine argument	1990	Frequency scale factor
ZSCALE	1	Subroutine argument	1991	Impedance scale factor
EPSLN	1	Subroutine argument	1992	Smallest non-zero matrix element

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Subroutine: CROLGR

Description: This subroutine calls the subroutines to read the source ordering, initial conditions, and run controls card groups and the end card. These card groups are signaled by header cards which contain asterisks in their first column. The names on these header cards and the expected order in which they appear are:

- * SOURCE ORDERING
- * INITIAL CONDITIONS
- * RUN CONTROLS
- * END

The source ordering, initial conditions, and run controls card groups need not appear in this order but must appear before the end card. In this order, however, they will process faster.

The source order card group must appear if more than one source is defined in the circuit description. Failure to do this will result in termination of program execution.

If the run controls card group does not appear, the following run controls are assumed:

- ISTOP = 1
- TO = 0.0
- TSTART = 0.0
- H = 10^{-4}
- HMIN = 0.0
- METHOD = 1
- IHWAX = 1
- MAXPTS = 100
- INCRIT = 3

A blank header card will be assumed to be the end card. If the end card does not appear, the end of file card is assumed to be the end card.

An unrecognizable header card will result in termination of program execution.

Usage: This subroutine is called by the main program when the Gear option has been chosen, immediately after control is returned from CORNAP.

Routine Name: ICONSV

Function: Read cards containing the names and initial values of the components of the initial state vector. Print all nonzero conditions used by the program and/or any diagnostic error messages generated.

Setting: At execution time the '* INITIAL CONDITIONS' card has just been processed, and the routine expects to read and process cards containing the initial circuit conditions, stopping at the first card read which contains an * in column 1.

Inputs:

Initial condition cards

NSTV - The number of state variables CORNAP obtains

NPSTV - The number of excess state variables*

NAMSV - Names of CORNAP obtained state variables

NAMLC - Names of excess state variables

Outputs: Error diagnostics and/or state vector of initial conditions. Components improperly formatted are set to zero. Excess components are ignored.

*An excess state variable is a capacitor voltage or inductor current whose value is dependent upon other capacitor voltages or inductor currents.

Subroutine: SRCORD

Function: This subroutine reads the source order cards to determine a correspondence between the way the sources are ordered in the state variable equations and the order in which they are computed and stored by USEFCN.

Inputs:

NIN - The number of independent sources defined in the circuit description.

NAMES - A vector containing the user names of the independent sources in the same order they appear in the state variable equations.

Outputs:

NORDER - A vector whose Ith element is the position in the Vector U(I) of USEFCN which corresponds to the position of the Ith source in the state variable equation as ordered by the circuit translation routines.

NUFFER - A vector containing the next control card.

Error messages are printed and the program terminated for the following conditions:

- (1) An input is specified more than once.
- (2) An input specified is not found in "NAMES."
- (3) Not all inputs defined in the circuit description are specified on the source order cards.

Additional inputs: This subroutine will read the source order cards. These cards contain the names of the sources, one per card in free format, in the order they are defined in USEFCN. All cards read are

assumed to be source order cards, unless they contain an asterisk in column 1. When a card containing an asterisk in column 1 is read, it is stored in NUFFER and the source order cards are assumed to be complete.

Subroutine: RUNCTL

Description. This subroutine reads the run controls card group for the Gear mode. The completion of this card group is signaled by a card with an asterisk in column 1 or the end of file card.

Input Parameters: The run controls cards must be one of the following:

INITIAL TIME = A
STEP RESPONSE
IMPULSE RESPONSE
PRINT START = A
STOP TIME = A
POINTS PRINTED = N
EPS = A
ADAMS INTEGRATION
HMAX = A
HMIN = A
HINIT = A

Here A is a floating point number and N is an integer. A description of the information conveyed by such cards is contained in the user's manual for STICAP.

Output Parameters:

TO - Initial value of time

TSTART - Time at which output printings is to begin

ISTOP - Stop condition

ISTOP = 1 Maximum number of points printed

ISTOP = 2 Maximum value of time

MAXPTS - Maximum number of points printed

TSTOP - Maximum value of time

H - Initial step size

EPS - Error control

METHOD - Integration method

METHOD = 0 Adams method

METHOD = 1 Stiff integration method

INCRMT - Number of integration steps between output printings

MAXDER - Maximum derivative used by the integration method.

MAXDER is set to 7 for METHOD = 1 and 8 for METHOD = 0.

HMAX - Maximum integration step size allowed

HMIN - Minimum integration step size allowed

NPUT - Type of input

NPUT = 0 User defined

NPUT = 1 Impulse

NPUT = 2 Step

NUFFER - Vector containing the next header card. If an end of
file card was encountered, NUFFER(2) is set to 1HE
which will be interpreted as the end card.

Default Options: Unless specified by the run controls cards, the
following values are assumed.

ISTOP = 1

TO = 0.0

TSTART = TO
NPUT = 0
HMAX = (TSTOP-TO)/10 if TSTOP was specified.
HMIN = 0.0
HMAX = 0.0 (interpreted as unbounded)
MAXPTS = 100
EPS = 10^{-4}
METHOD = 1
INCRMT = 3
H = 10^{-4}
MAXDER = 7

Machine Dependence: This subroutine itself is machine independent but calls the machine dependent routines APACK and ATODEC.

Subroutines Required: APACK, ATODEC

Program: CROLMY

This subroutine calls the subroutines to read the source definition, initial conditions, and run controls card groups and the END card. These card groups are signaled by header cards which contain asterisks in their first columns. The names on these header cards are

- * SOURCE DEFINITIONS
- * INITIAL CONDITIONS
- * RUN CONTROLS
- * END

The END card must be the last card appearing. If it is not present, the end of file card is interpreted as the END card.

The run controls card group must be present. The source definition card group must be present unless the impulse or step response was specified in the run controls card group. If not execution will be terminated.

Subroutine: SRCINT

Description:

This subroutine interprets the source definition card group for the matrix mode. This card group is signaled by a card with an asterisk in column 1, followed by the words "SOURCE DEFINITIONS" in free format. The subroutine assumes all cards following this card up to the next card with an asterisk in column 1 are source definition cards. These cards are of the form

$$\text{NAME} = f_1(t) + f_2(t) + \dots + f_N(t)$$

where NAME is the name of a source defined in the circuit description, and each $f_i(t)$ is one of the following

A

A*IMP

A*SIN(B*T)

A*SIN(B*T+C)

A*COS(B*T)

A*COS(B*T+C)

where A, B, and C are floating point numbers in free format. The asterisks indicating multiplication need not appear. A source description may be continued on additional cards by placing a dollar sign (\$) in column 1 of the cards on which the description is

continued as long as each $f_1(t)$ is completely described on only 1 card and the total number of functions describing a given source is no more than 20. A sinusoidal or cosinusoidal function with a phase angle is broken into a sinusoidal plus a cosinusoidal function with zero phase angle thus counting as two functions.

This subroutine will terminate program execution if an error occurs in a source description or if not all the sources defined in the circuit description are described in the source definition cards.

Additional Input Parameters

- NIN - The number of circuit inputs.
- NAMIN - A vector containing the names of the circuit inputs in the order they appear in the state variable equations.

Output Parameters

- NORDER - A vector relating the order the sources are defined in the state variable equations with the order they are defined in NSORCE, ISORCE, and SOURCE.
- ISORCE - The number of functions describing source I is ISORCE(K) where $K = \text{NORDER}(I)$
- NSORCE - The function type for the Jth function describing the Ith source is prescribed by NSORCE(K,J) where $K = \text{NORDER}(I)$, and
- | | |
|-----------------|---------|
| NSORCE(K,J) = 1 | step |
| NSORCE(K,J) = 2 | sine |
| NSORCE(K,J) = 3 | cosine |
| NSORCE(K,J) = 4 | impulse |

SOURCE - The magnitude of the Jth function of the Ith source is contained in SOURCE(K,J,1), where K = NORDER(I) and the angular velocity (for sine and cosine functions) is SOURCE(K,J,2).

Machine Dependence - This subroutine is machine independent but calls the machine dependent subroutines APACK and ATODEC.

Subroutine: RUNCRL

This subroutine reads in the run controls card group for the matrix method.

The completion of this card group is signaled by a card with an asterisk in column 1 or the end of file card.

Input Parameters

The run controls cards read must be one of the following

INITIAL TIME = A
STEP RESPONSE
IMPULSE RESPONSE
PRINT START = A
STOP TIME = A
POINTS PRINTED = N
PLOT INCREMENT = A

where A is a floating point number and N is an integer.

Output Parameters

TO Initial value of time
TSTART Time at which output printings is to begin.
NPTS Number of output points to be printed.
INPUT Type of input
DT Time between output printings.

Default Options

If no initial time is specified, it is assumed to be zero.

If no print start time is specified, it is assumed to be the initial time.

If a stop time but no plot increment or points printed is specified or if only the plot increment is specified, 100 time points will be printed.

Machine Dependence

This subroutine is machine independent but calls the machine dependent routines APACK and ATODEC.

Program: CROLGM

This program calls COMPAK to rearrange the variables in blank common from the form supplied by CORNAP to the form required by GEAR and MATRIX. One of the subroutines CROLGR or CROLMX is then called to read the card groups peculiar to the GEAR or the MATRIX mode.

Program INTGR

This subroutine performs the integration of the state variable equations.

The upper limit of the integration may be given by specifying the maximum value of the independent variable time or by specifying the number of points printed.

If the maximum value of time is the stop condition, interpolation is used to determine the values of the state variables at the stop time.

The following error conditions will terminate program execution

1. Requested error is smaller than can be handled for this problem.
2. Corrector convergence could not be achieved with H greater than HMIN.
3. The requested error could not be achieved with H equal to hMIN.

Input Parameters

NSTV	The number of state variable equations.
TO	The lower integration limit.
XO	A vector containing the values of the state variables measured at time TO.
TSTART	The time at which printings is begun.
TSTOP	The upper integration limit
LAXPTS	The maximum number of time points to be printed.
ISTOP	The stop condition

ISTOP = 1 Maximum time points printed.

ISTOP = 2 Maximum value of time.

NO The number of circuit outputs

NAMOUT A vector containing the names of the circuit outputs

H The initial integration step size.

HMIN The minimum integration step size.

HMAX The maximum integration step size permitted.

INCRMT The number of integration steps between printings.

EPS The euclidean norm of a vector whose i th element is the single step error of the i th state variable divided by the maximum value of the i th state variable previously encountered must be less than this value.

METHOD The integration method used.

METHOD = 0 An Adams predictor corrector method.

METHOD = 1 A multistep method suitable for stiff systems.

MAXDER The maximum derivative that should be used in the method. It must be less than 8 for Adams and 7 for stiff methods.

Subroutine: DATPNT

This subroutine prints the output for the Gear mode. It performs the calculation

$$\underline{Y}(t) = \underline{C} \underline{x}(t) + \underline{D} \underline{u}(t)$$

where $\underline{Y}(t)$ is the output vector at time T . $\underline{x}(t)$ is a vector containing the values of the state variables at time T , and $\underline{u}(t)$ is a vector containing the values of the inputs at time T .

Input Parameters:

T - present value of time.
x - a 8 x 30 matrix. x(1,I) contains the value of the Ith state variable measured at time T.
c,D - constant matrices of the output equation.
NO - number of circuit outputs.
NSTV - number of state variable equations.
NIN - number of inputs defined in the circuit description.
NORDER- a vector telling the order the inputs are defined in USEFCN.

Output Parameters:

This subroutine prints the value of time and the values of the outputs measured at this time.

Subroutine Required: USEFCN

Machine Dependency: None

Subroutine: DIFFUN(T,X,DX)

This subroutine evaluates the derivatives of the dependent variables X, with respect to the independent variable T, using the matrix equation

$$\dot{X} = AX + BU$$

Input Parameters

NSTV The number of state variable equations.
NIN The number of circuit inputs
A System matrix
B Input matrix

T Present value of time

X X(1,I) contains the value of the Ith state
 variable measured at time T.

U U(I) is the value of the NORDER(I)TH input
 measured at time T

NORDER NORDER(I) is where the Kth source is located

Output Parameters

DY DY is a vector containing \dot{X} measured at the
 present values of X and T.

Subroutines Required

USEFCN

Soubroutine: PEDERV

Description: This subroutine supplies DIFSUB with the Jacobian
matrix of the state variable equation:

$$\dot{\underline{X}} = \underline{A} \underline{x} + \underline{B} \underline{u}$$

which is, in fact, the A matrix.

Input Parameters:

M - number of state variable equations.

A - system matrix stored in core with the first column of A
 in the 1st 30 storage locations, the second column of A
 in the next 30 storage locations and so forth.

Output Parameters:

PW - Jacobian matrix. This is the system matrix, A, with
 each column stored in core starting in the core location
 following the end of the previous column.

Subroutines Required: None

Machine Dependency: None

Example of Usage:

Given the system matrix

$$A = \begin{bmatrix} 1 & 4 & 7 \\ 2 & 5 & 3 \\ 3 & 6 & 9 \end{bmatrix}$$

A is supplied as

1,2,3,27 blanks,4,5,6,27 blanks,7,8,9,810 blanks,b

PW is outputted as

1,2,3,4,5,6,7,8,9,891 blanks,b

Subroutine: MATINV(A,N,JSING)

Purpose

Invert a matrix

Description of parameters

A - Input matrix, destroyed in computation and replaced by resultant inverse.

N - Order of matrix A

JSING - Indicates whether A is singular; one if not, minus one if it is.

Remarks

Matrix A must be a general matrix

Subroutines and function subprograms required: None

Method

The standard Gauss-Jordan method is used. The determinant is also calculated. A determinant of zero indicates that the matrix is singular.

Subroutine: DECOMP(N,PSAVE,J1)

Function:

This routine performs an LU decomposition of the NxN matrix found in PSAVE. PSAVE is replaced by the matrices U and L, with U the main diagonal and above, L below the main diagonal. All main diagonal elements of L are unity, so need not be stored.

If the PSAVE matrix is nonsingular, the routine returns a value of +1 for J1; if singular, J1 is set to -1.

The routines DECOMP and SOLVE are lifted directly from Forsythe and Moeller, Numerical Linear Algebra, p. 68 and are well discussed in this reference.

Subroutine: SOLVE(N,PSAVE,SAVE,N5)

Function.

This routine employs the LU decomposition of DECOMP stored in PSAVE to solve for \bar{X} the N equations

$$(LU)\bar{X} = \bar{b},$$

with various \bar{b} supplied by Gear's subroutine. The result is stored in SAVE. N5 is a parameter used for addressing the proper locations in SAVE in which \bar{X} is to be stored

For large systems (N greater than 5) the LU decomposition is faster than the (MATINV) matrix inversion previously used by Gear. If desired, minor alterations to STICAP will restore the Gear routine to its previous form. These modifications are stated in Gear's ACM DIFSUB description (see STICAP theory manual for the complete reference).

Program: MATRIX

This overlay solves the system of state variable equations by performing a spectral decomposition of the transition matrix then integrating in closed form. This technique is described in the paper:

"An Efficient Matrix Algorithm for the Solution of Linear Systems with Widely Separated Eigenmodes" ... by E. Young, J. Heinbockel, and M. Ransom, Proceedings of the 4th Asilomar Conference on Circuit and System Theory, Pacific Grove, Calif., 1970.

Inputs to this program are described in the comments of the three secondary overlays and in the programmer's manual.

This first secondary overlay determines the eigenvalues of the system matrix, determines the inverse of the Vandermonde matrix given by these eigenvalues, and finally determines a sparse matrix code for the system matrix.

CALL OVERLAY(3HCOM,4,1,0)

The next secondary overlay forms the transition matrix times a vector

$$\text{EXP}(A*T)*V$$

as a finite series for the initial condition vector and for each driving vector. This expression is then integrated in close form and the symbolic solution evaluated at each time point at which output is requested.

CALL OVERLAY(3HCOM,4,2,0)

This final secondary overlay prints the tabulated results, and if requested, the close form solution.

CALL OVERLAY(3HCOM,4,3,0)

Program: SETSVE

This program calls EIGNV to determine the eigenvalues of the system matrix and VANINV to form the inverse of the Vandermonde matrix given by these eigenvalues. Finally this program determines a sparse matrix code for the system matrix.

Input Parameters:

- A - The system matrix.
- N - The number of state equations.
- EPS - The value below which an element is felt to be zero.

Output Parameters:

- C - A complex matrix containing the inverse of the Jacobian matrix.
- KNZ - The number of non-zero elements in the system matrix.
- NZR,NZC - Vectors containing a sparse matrix code for the system matrix. The Ith non-zero element of the system matrix is located in row NZR(I) and column NZC(I).
- ANZ - A vector which replaces A in core which contains only the non-zero elements of A.

Subroutine: VANINV

This subroutine forms the inverse of the Vandermonde matrix given by the eigenvalues of the system matrix. The Vandermonde matrix of the n elements $\lambda_1, \lambda_2, \dots, \lambda_n$ is the n x n matrix

$$V = \begin{bmatrix} 1 & 1 & \dots & 1 \\ \lambda_1 & \lambda_2 & \dots & \lambda_n \\ \lambda_1^2 & \lambda_2^2 & \dots & \lambda_n^2 \\ \vdots & \vdots & & \vdots \\ \lambda_1^{n-1} & \lambda_2^{n-1} & \dots & \lambda_n^{n-1} \end{bmatrix}$$

The inverse of the Vandermonde matrix is formed in this subroutine using a technique due to Kaufman.¹

$$\text{Let } [C_{ij}] = [V_{ij}]^{-1}$$

$$\text{then } C_{ij} = \frac{\sum_{k=0}^{n-j} a_k \lambda_i^{n-j-k}}{n \prod_{\substack{k=1 \\ k \neq i}}^n (\lambda_i - \lambda_k)}$$

where the a_k are the coefficients of the characteristic equation

$$\begin{aligned} P(\lambda) &= \prod_{i=1}^n (\lambda - \lambda_i) \\ &= a_0 \lambda^n + a_1 \lambda^{n-1} + \dots + a_{n-1} \lambda + a_n \end{aligned}$$

In this subroutine the a_k are formed from the recursive relationship

$$\begin{aligned} a_0 &= 1 \\ a_k &= -\frac{1}{k} \sum_{m=1}^k a_{k-m} T_m \end{aligned}$$

$$\text{where } T_m = \sum_{p=1}^n \lambda_p^m$$

We note finally that if $\lambda_i = \lambda_j^*$, then $C_{ik} = C_{jk}^*$.

Input Parameters:

N - The number of state variable equations.

ROOT - A complex vector whose elements are the eigenvalues of the system matrix.

Output Parameters:

C - A complex matrix containing the Vandermonde inverse.

-
1. Kaufman, I., "Evaluation of an Analytical Function of a Companion Matrix with Distinct Eigenvalues," 1969 Proceedings of IEEE Letters, pp. 1180-1181.

Subroutine: HESEN(A,M)

Supplier: NASA Routine*

This subroutine reduces a matrix through similarity transformations to an upper-Hessenburg matrix. An upper-Hessenburg matrix is one in which all elements below the subdiagonal are zero.

This subroutine is the NASA Langley Research Center routine HESSEN.

Input Parameters:

M - The order of the matrix being reduced.

A - The matrix being reduced. This matrix is destroyed in the computations.

Output Parameters:

A - The reduced matrix which has replaced in input matrix in core.

Subroutine: QRTRAN(A,N,R,SIG,D)

Supplier: NASA Routine

This subroutine performs one double Q-R transformation on the matrix A.

$$\begin{aligned}\text{Let } A &= Q_1 R_1 \\ A_1 &= R_1 Q_1 = Q_1^{-1} A Q_1 = Q_2 R_2 \\ A_2 &= R_2 Q_2 = Q_2^{-1} A_1 Q_2 = Q_2^{-1} Q_1^{-1} A Q_1 Q_2\end{aligned}$$

where

Q_1 and Q_2 are unitary matrices

and

R_1 and R_2 are right triangular matrices.

*Not contractor supplied.

A is then replaced in core by A_2 .

This subroutine is the NASA Langley Research Center routine QR2.

Subroutine: EIGNV

Supplier: NASA Routine

This subroutine finds the eigenvalues of a real, non-symmetric matrix using the double Q-R transformation.

This routine was slightly modified from the NASA Langley Research Center routine REIS.

Input Parameters:

KJN - The number of state variable equations.

AO - The matrix whose eigenvalues are being determined.

Output Parameters:

ROOT - A complex vector containing the eigenvalues of the input matrix with real eigenvalues stored first and each complex eigenvalue followed immediately by its complex conjugate.

Program: INTCL

This overlay calls the various integration routines to perform the closed form integration for each function making up each circuit input. The close form solution is then evaluated at each time point requested.

Input Parameters:

N The number of state variables.

NPTS The number of time points to be printed.

TSTART	The time at which output printings are to be begun.
DT	The time increment between output printings.
A,B,C,D	Matrices in the input and output equations.
NIN	The number of circuit inputs.
NO	The number of circuit outputs.
XO	A vector containing the initial conditions of the state variables.
INZ	Tells whether all initial conditions are zero. INZ = 1 if there are any non-zero initial conditions.
NZR,NZC,KNZ	Variables containing a sparse matrix code for the system matrix.
IPRINT	Tells whether or not the closed form solution is to be printed.
NORDER	A vector relating the way the sources are defined in ISORCE, NSORCE and SOURCE with the way they are defined in the output and state variable equations.
ISORCE,SOURCE, NSORCE	Matrices containing information about sources. See SCEVAL for description.
NSC	The number of different angular velocities exciting the circuit.

Output Parameters:

OUT	A matrix containing the outputs evaluated at each time point.
-----	---

Subroutines required in this overlay:

STEADY, INIVAL, STORE, TRANS, SCEVAL

EAT, SVESTP, SVESIN, SVECOS, IPROD

Subroutine: SVECOS(AMPL,OMEGA)

This subroutine determines the response of the system due to cosine driving functions.

Input Parameters:

N The number of state variable equations.
NR The number of real eigenvalues of the system matrix.
AMPL The amplitude of the driving function.
OMEGA The angular velocity of the cosine driving function.
P A NXN complex matrix whose columns are the V'S of the
 expression

$$\begin{aligned} \text{EXP}(A*T)*R &= V1*\text{EXP}(LAMBDA(1)*T) \\ &+ V2*\text{EXP}(LAMBDA(2)*T) \\ &+ \dots \\ &+ VN*\text{EXP}(LAMBDA(N)*T) \end{aligned}$$

where

A is the system matrix.

R is the driving vector.

LAMBDA(I) is the Ith eigenvalue of A.

and

T is time.

TO The time at which the initial conditions of the system
 were measured.

ROOT A complex vector of length N containing the eigenvalues
 of the system matrix.

Output Parameters:

CFCOS A NXN matrix containing the transient part of the solution.

STYCOS A NX2 rectangular matrix containing the steady-state portion of the solution.

Subroutine: SVESIN(AMPL,OMEGA)

This subroutine performs the integration of $\text{EXP}(A(T-\text{TAU})) * R * \text{SIN}(\text{OMEGA} * X)$ from T_0 to t where A is a matrix and R is a vector.

Input Parameters:

N The number of state variables.

OMEGA The angular velocity of the driving function.

R A vector of N length associated with the driving function.

TO The initial value of time.

ROOT A vector containing the circuit eigenvalues.

P A matrix containing $\text{EXP}(A * T) * R$.

NR The number of real eigenvalues.

AMPL The amplitude of the driving function.

Output Parameters:

STYSIN A NX2 rectangular matrix containing the steady state solution to the integration.

COEF A NXN matrix containing the transient part of the solution.

Subroutine: SVESTP(AMPL)

This subroutine determines the responses due to step driving sources.

Input Parameters:

- N . The number of state variable equations.
- NR The number of real eigenvalues of the system matrix.
- ROOT A complex vector of length N containing the eigenvalues of the system matrix.
- TO The time at which the initial values of the state variables are measured.
- P A NXN complex matrix whose columns are the V'S of the expression

$$\begin{aligned} \text{EXP}(A*T)*R &= V1*\text{EXP}(\text{ROOT}(1)*T) \\ &+ \dots \\ &+ VN*\text{EXP}(\text{ROOT}(N)*T) \end{aligned}$$

where

A is the system matrix

R is the driving vector

ROOT(I) is the Ith eigenvalue of A

and

T is time.

Where complex conjugates occur, the complex conjugate V is omitted and the remaining V'S stored consecutively in P

AMPL The amplitude of the step function.

Output Parameters:

- COEF A NXN matrix containing coefficients corresponding to the transient part of the response.
- STP A vector of length N containing the steady state dc part of the response.

Subroutine: INIVAL(AMPL)

This subroutine is called to determine the solution corresponding to the initial conditions or for impulse driving functions.

Input Parameters:

N Number of state variables.
ROOT Vector containing the eigenvalues of the system matrix.
NR Number of real eigenvalues.
P A matrix containing $\text{EXP}(A(T-T_0))$ where A is the system matrix.
AMPL The amplitude of the impulse function.

Output Parameters:

COEFO A matrix containing the closed form solution.

Subroutine: MPROD(A,B,L,M,N)

This subroutine multiplies a matrix A into a matrix B and stores the product in B.

Input Parameters:

A A matrix dimensioned 10x30.
B A matrix dimensioned to have 30 rows but having a variable number of columns. This matrix is destroyed during the computation.
L Number of rows in matrix A.
M Number of columns in matrix A and rows in matrix B.
N Number of columns in matrix B.

Output Parameters:

B A matrix containing the product.

Subroutine: EAT(V)

This subroutine finds the expression

$$e^{\underline{A}t} \cdot \underline{V}$$

where \underline{A} is the system matrix and \underline{V} is a given vector. This expression is given as

$$e^{\underline{A}t} \cdot \underline{V} = \underline{V}_1 e^{\lambda_1 t} + \dots + \underline{V}_n e^{\lambda_n t}$$

Here

$$\underline{V}_1 = \sum_{k=1}^n C_{1k} \underline{A}^{k-1} \underline{V},$$

$\lambda_1, \lambda_2, \dots, \lambda_n$ are the eigenvalues of \underline{A} and $[C_{ij}]$ is the inverse of the Vandermonde matrix given by these eigenvalues. Since if $\lambda_j = \lambda_i^*$, $\underline{V}_j = \underline{V}_i^*$ complex conjugate \underline{V} 's are not given.

Input Parameters:

- | | |
|---------|--|
| N | The number of state variable equations. |
| NR | The number of real eigenvalues in the system matrix. |
| A | A vector containing the non-zero elements of the system matrix. |
| KNZ | The number of non-zero elements in the system matrix. |
| NZR,NZC | Two vectors containing a sparse matrix code for the system matrix. The Ith non-zero element in the system matrix is in row NZR(I) and column NZC(I). |
| C | A complex matrix containing the inverse of the Vandermonde matrix given by the eigenvalues of the system matrix. |

Output Parameters:

- | | |
|---|--|
| P | A matrix whose columns are the \underline{V} 's in the above expansion except that if \underline{V}_i is complex, $\text{Re}[\underline{V}_i]$ is stored |
|---|--|

in the Ith column of \underline{P} and $\text{Im}[\underline{V}_1]$ is stored in the I+1th column of \underline{P} . \underline{V}_{i+1} which is \underline{V}_1^* is then omitted.

Subroutine: SCEVAL

This subroutine evaluates the expression

$$D*U$$

where $D*U$ is the matrix expression in the output equation

$$Y = C*X + D*U$$

at each time point at which output is requested and adds the result to the contents of a matrix out which contains $C*X$ at each output time point.

Input Parameters:

TSTART The time at which output printings are to start.

DT The time increment between output printings.

NUM The number of output time points to be printed.

NIN The number of circuit inputs.

NO The number of circuit outputs.

NORDER A vector relating the way the sources are defined in ISORCE, NSORCE, and SOURCE with the way they are ordered in the state and output equations.

ISORCE A vector whose Ith element tells the number of function making up the NORDER(I) source.

NSORCE A matrix whose I,J element tells the type of the Jth function making up the NORDER(I) source.

NSORCE(I,J)=1 STEP

NSORCE(I,J)=2 SINE

NSORCE(I,J)=3 COSINE

NSORCE(I,J)=4 IMPULSE

SOURCE A 3-dimensional matrix whose I,J,1 element tells the magnitude of source NSORCE(I,J), and whose I,J,2 element tells (for sine and cosine functions) the angular velocity of the function.

D The matrix relating the outputs to the inputs in the output equation.

Output Parameters:

OUT A matrix containing the circuit outputs at the requested time points.

Subroutine: TRANS

This subroutine calculates the transient portion of the solution by substituting into the closed form expressions for the solutions.

Input Parameters:

N The number of state variable equations.

NUM The number of different time values at which the output is to be measured.

TSTART The time at which the first measurement is to be made.

COEF A matrix containing the coefficients of the transient portion of the solution.

ROOT A complex vector of N length containing the eigenvalues of the system matrix.

NR The number of real eigenvalues of the system matrix.

DT The interval between time samples.

Output Parameters:

OUT A matrix containing the outputs at the requested time points.

The values of the outputs are given at time

$T_{START} + K \cdot DT$, $K=0,1,\dots,NUM-1$

Additional Subroutines Required: None

Subroutine: STEADY(NTYPE,OMEGA)

This subroutine calculates the steady state portion of the solution by substituting into the closed form solution. This calculation is made at time

$T_0 + K \cdot DT$ $K = 0,1,\dots,NUM-1$

where T_0 is the time at which printing is to begin,

NUM is the number of time points at which printing is
to be made,

and DT is the printing increment.

Input Parameters:

N The number of state equations.

NUM The number of different time points at which printing
is to be made.

TSTART The time at which the first printing is to be made.

NTYPE The type of driving function.

NTYPE = 1 Sine or cosine driving function.

NTYPE = 2 Step function.

STYSC A matrix containing the closed form steady state portion
of the solution for sine and cosine driving functions..

STP A vector containing the closed form steady state portion
of the solution for step functions.

Output Parameters:

OUT A matrix containing the circuit outputs at the requested
time points.

Subroutine: STORE(OMEGA)

This subroutine stores the steady-state response for sine and cosine driving functions. The subroutine will combine responses with the same angular velocities.

Input Parameters:

- NO The number of circuit outputs.
- STYSC A matrix whose I,1 element contains the steady-state sine response for the Ith output and whose I,2 element contains the corresponding cosine response.
- OMEGA The angular velocity of the driving function.

Output Parameters:

- NSC The number of different angular velocities encountered so far.
- ANGVEL A vector whose Ith component contains the angular velocity of the Ith sine or cosine steady-state response.
- SICO A matrix whose 1,J,K element contains the steady-state response of the Kth output with the Jth angular velocity and whose 2,J,K element contains the corresponding cosine response.

Program: OUTPNT

This overlay is used for printing output information for the matrix mode. Tabulated results are always printed. Close-form solutions are printed only if requested.

Input Parameters:

- IPRINT Tells whether or not the closed form solution is to be printed.

See CLSOL and TABPNT for additional input parameters.

Subroutine: CLSOL

This subroutine prints the closed form solution for the requested outputs.

Input Parameters:

NO The number of circuit outputs.

NAMOUT A vector containing the names of the circuit outputs.

NSC The number of sine and cosine inputs.

ANGVEL A vector whose Ith element is the angular velocity of the Ith sine or cosine input.

SICO A matrix whose 1,J,I element tells the steady-state sinusoidal response of the Ith output due to the Jth sine or cosine input, and whose 2,J,I element tells the corresponding steady state cosinusoidal response.

STP A vector whose Ith element gives the steady-state step response for the Ith output.

N The number of state variables.

NR The number of real eigenvalues.

ROOT A complex vector containing the eigenvalues of the system matrix with the real eigenvalues stored first and each complex eigenvalue followed immediately by its complex conjugate.

COEF A matrix containing the transient solution. The transient solution for the Ith output is

$$\text{COEF}(I,J) * \text{EXP}[\text{ROOT}(J) * T]$$

for ROOT(J) a real eigenvalue. When ROOT(J) is a complex

eigenvalue the solution corresponding to $\text{ROOT}(J)$ is combined with the solution corresponding to the complex conjugate of $\text{ROOT}(J)$ to give:

$$\begin{aligned} & \text{EXP}[\text{RE}(\text{ROOT}(J)) * T] \\ & * [\text{COEF}(I, J) * \text{SIN}(\text{IM}(\text{ROOT}(J)) * T) \\ & + \text{COEF}(I, J+1) * \text{COS}(\text{IM}(\text{ROOT}(J)) * T)] \end{aligned}$$

where $\text{RE}(\text{ROOT}(J))$ is the real part of $\text{ROOT}(J)$ and $\text{IM}(\text{ROOT}(J))$ is the imaginary part of $\text{ROOT}(J)$.

Subroutine: TABPNT

This subroutine prints the tabulated results for the matrix mode.

Input Parameters:

TSTART Time at which printing is to begin.
DT Time increment between output printings.
NPTS The number of time points to be printed.
OUT A matrix whose I,J element is the Jth output measured at the Ith time value.
NO The number of circuit outputs.
NAMOUT A vector containing the names of the circuit outputs.

Subroutine: ATAN3(X,Y)

This subroutine combines a sine function with a cosine function where both functions have equal angular velocities but no phase angle into a sine function with a phase angle. The phase angle is chosen to be positive or negative such as to make the magnitude of the phase angle less than $\pi/2$.

Input Parameters:

X The coefficient of the sine function.

Y The coefficient of the cosine function.

Output Parameters:

X The coefficient of the resulting sine function.

Y The phase angle in radians of the resulting sine function.

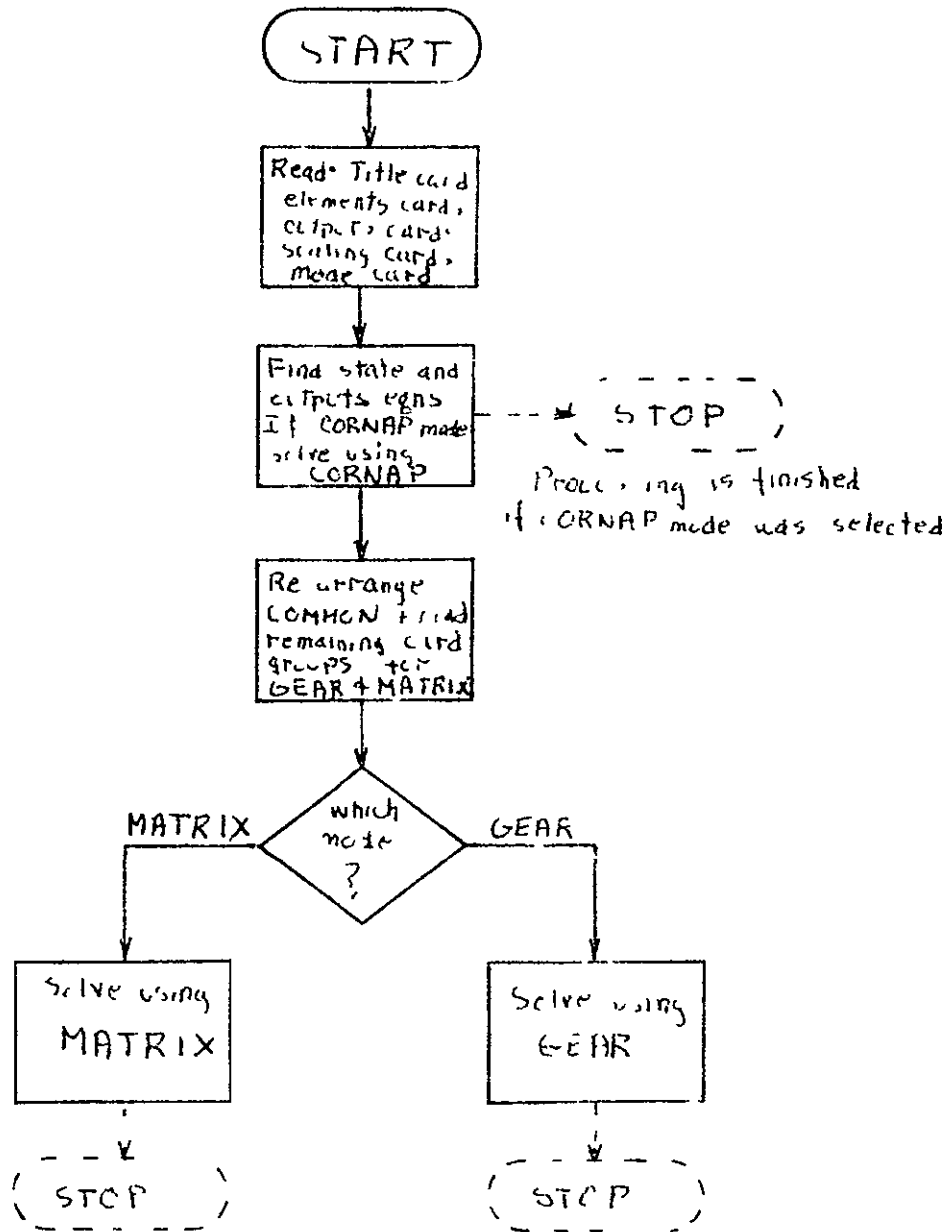
Note: Both input parameters are destroyed in the computations.

CHAPTER VI

FLOWCHARTS

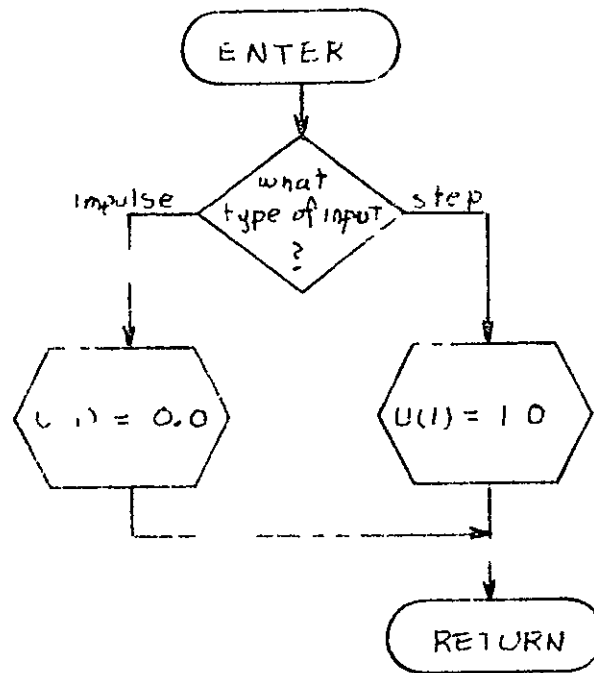
This chapter contains the flowcharts for all contractor supplied subroutines.

STICAP



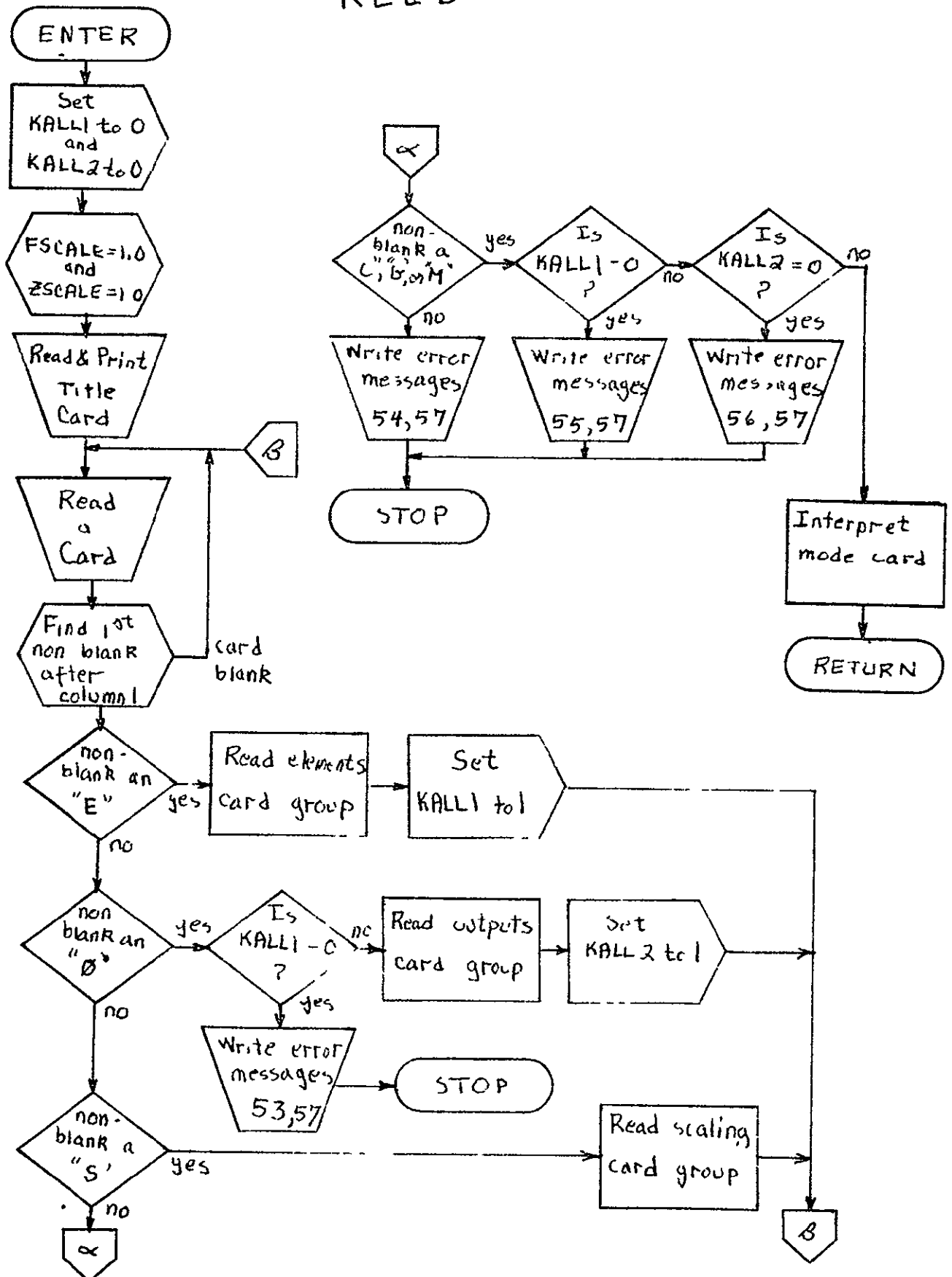
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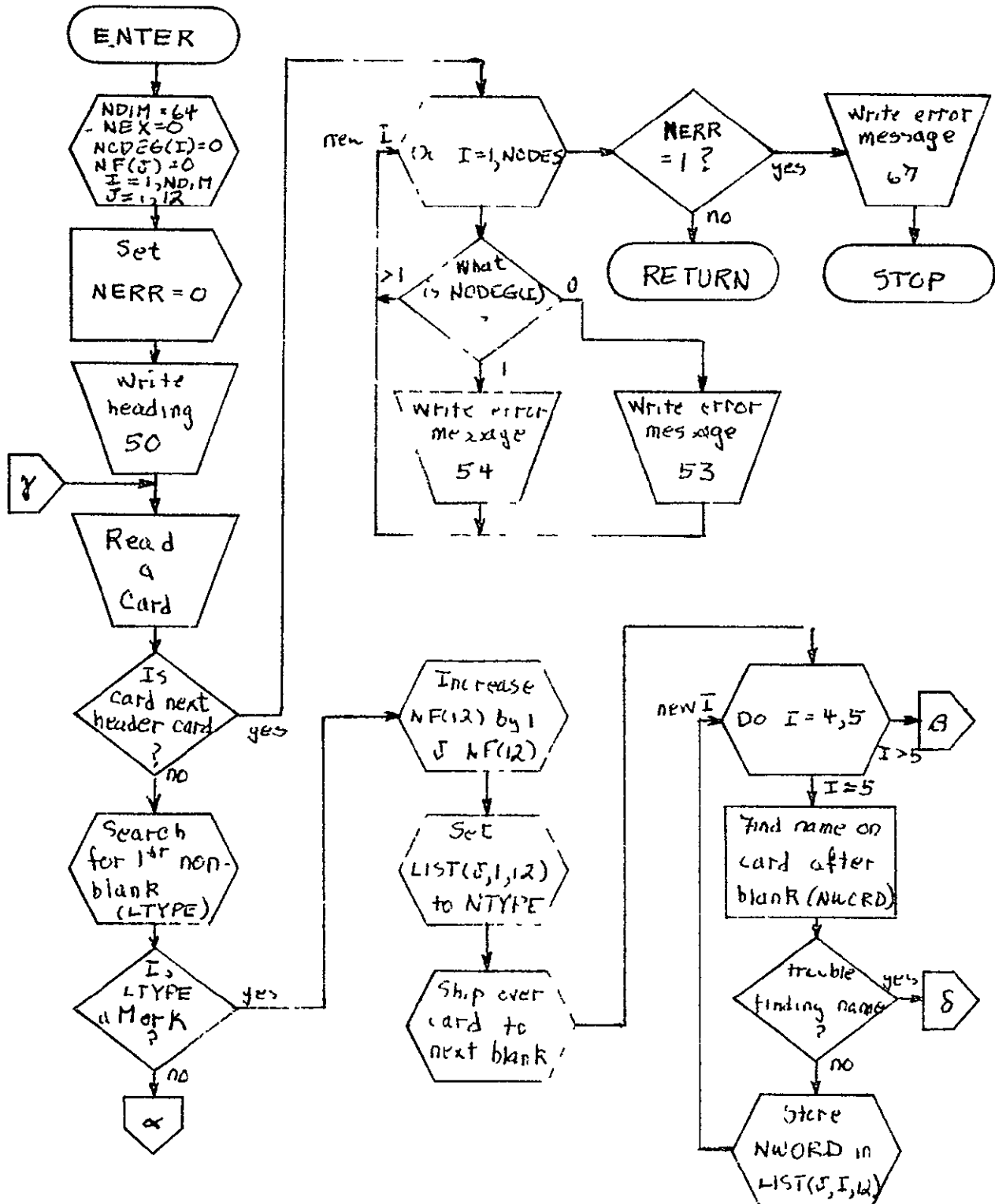


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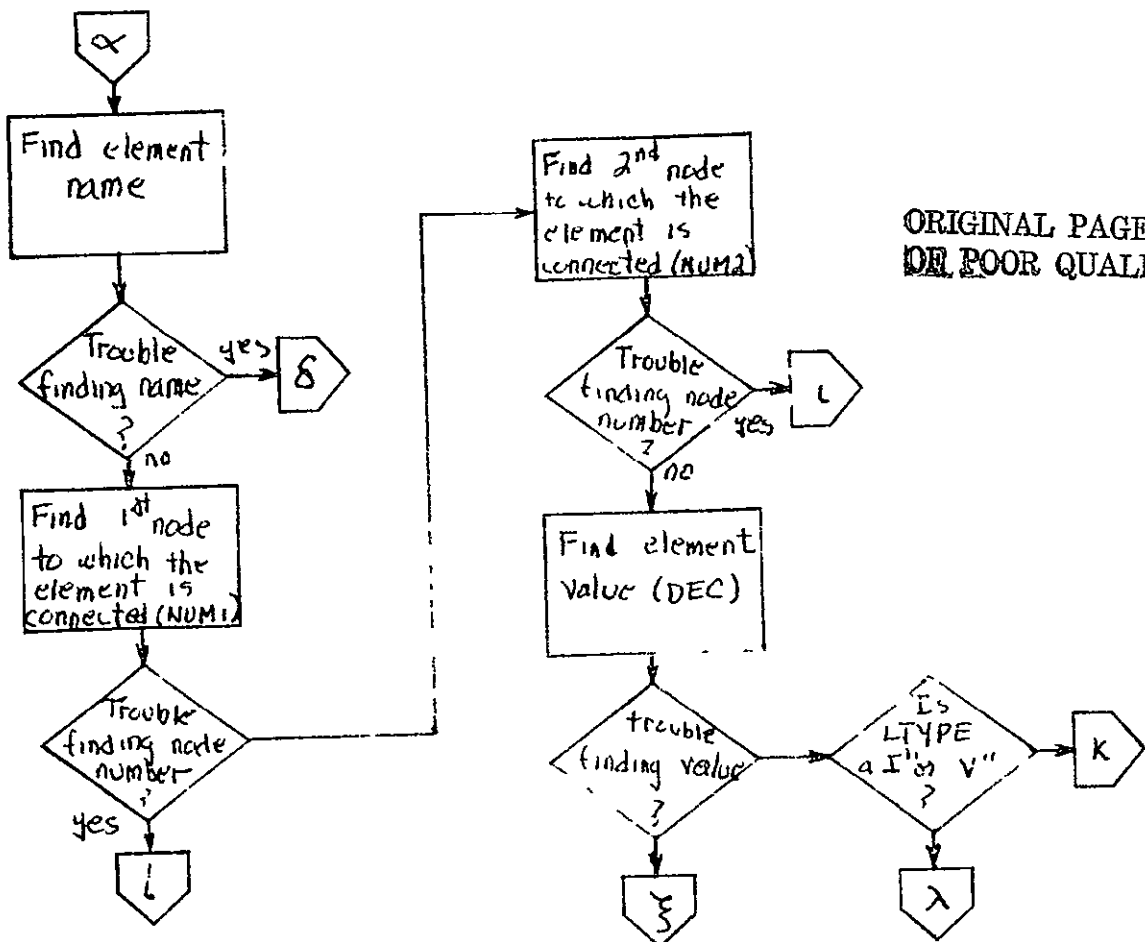
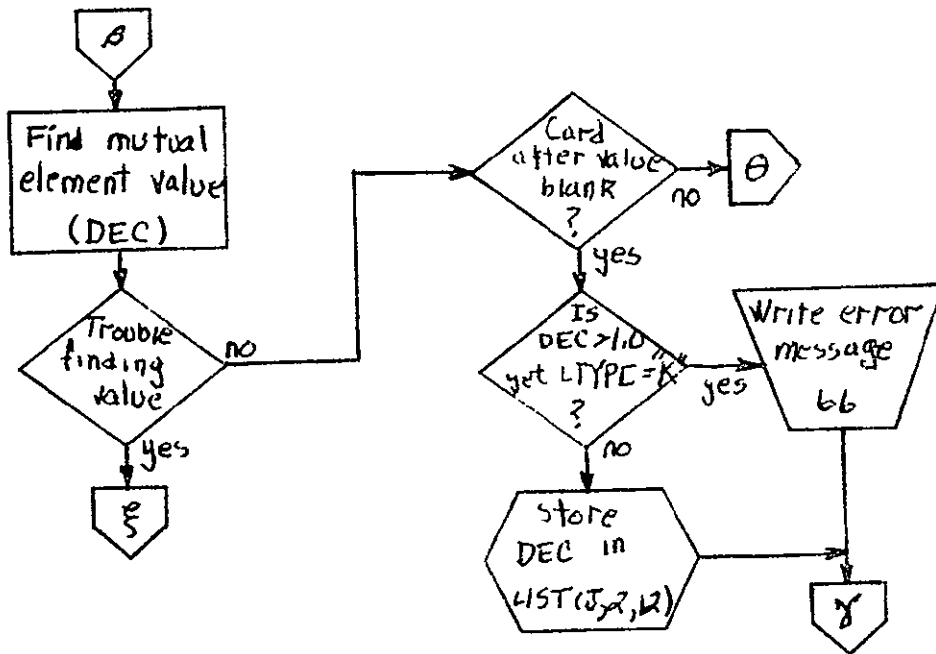
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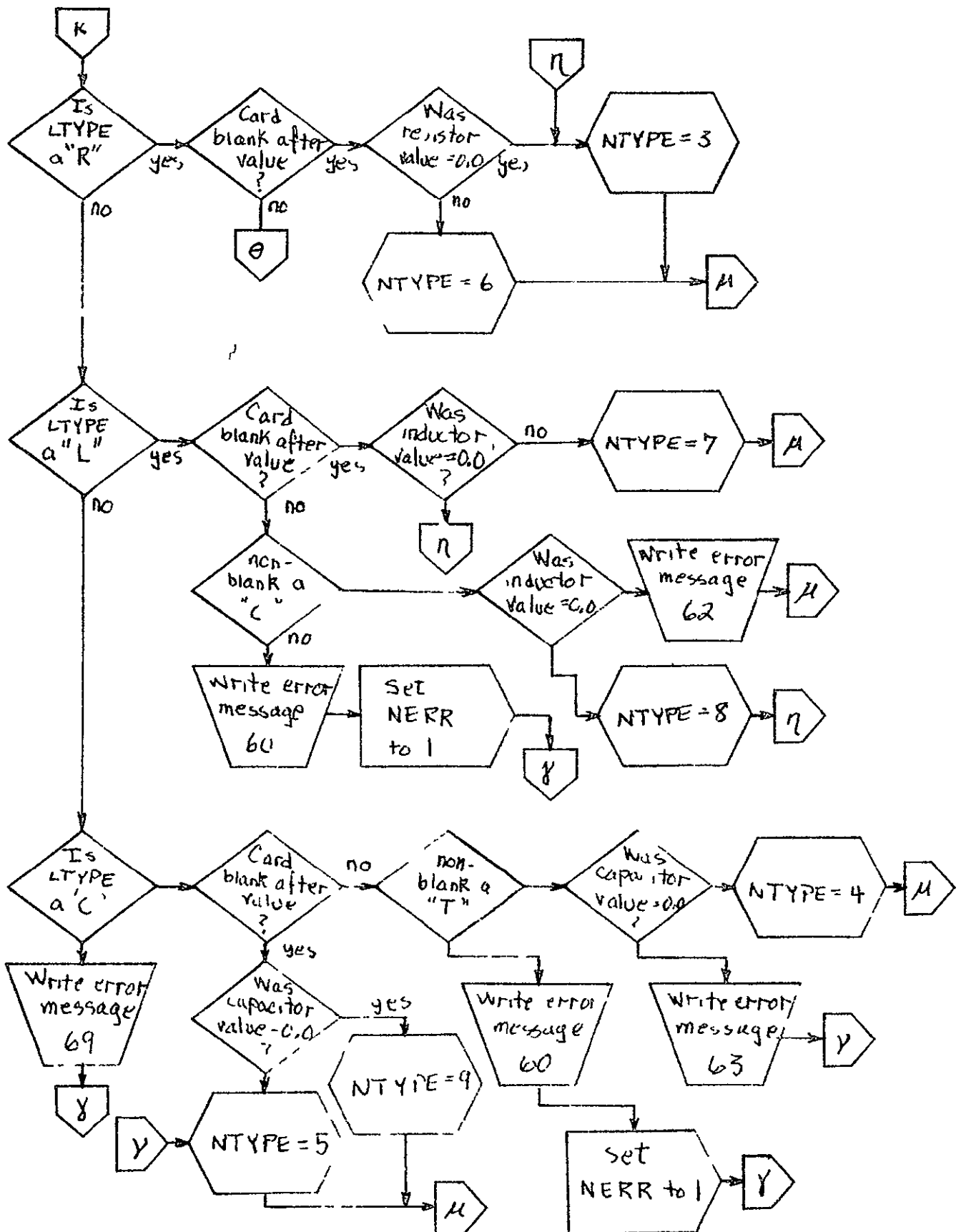
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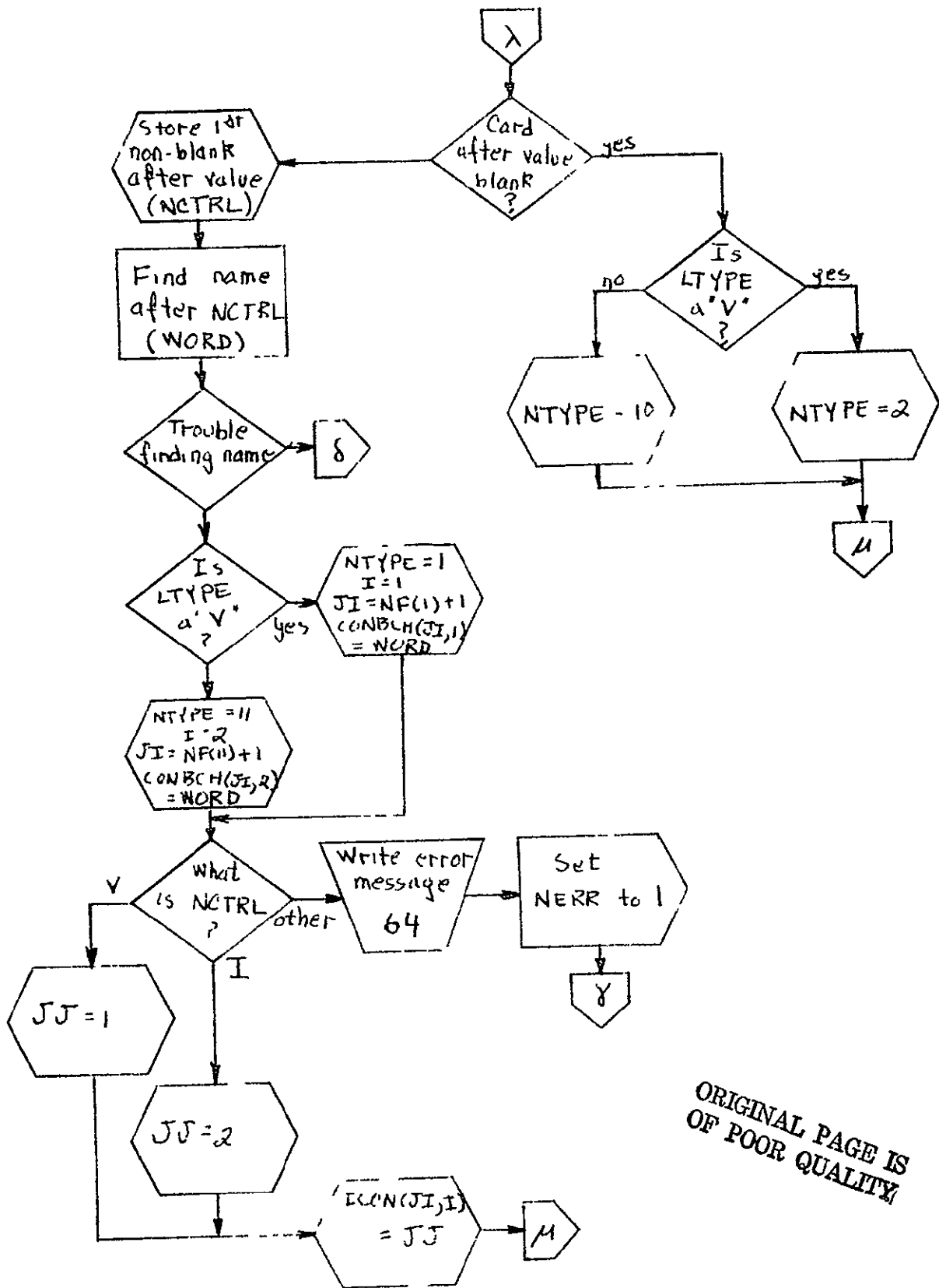


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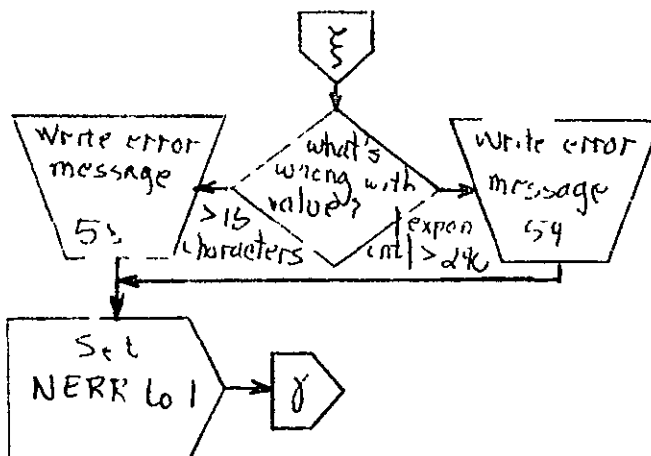
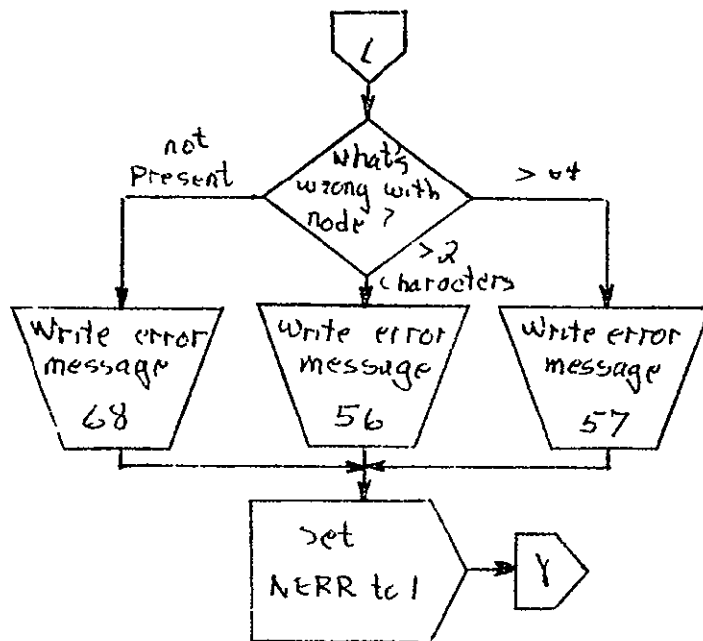
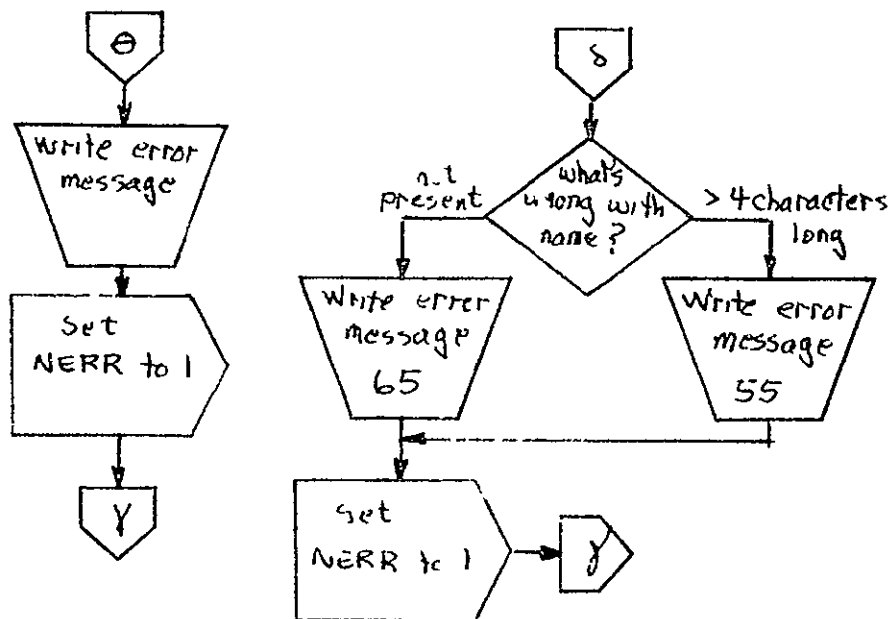
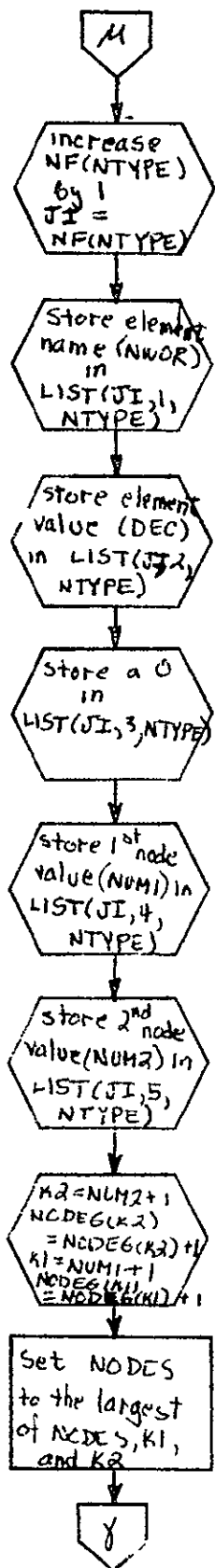


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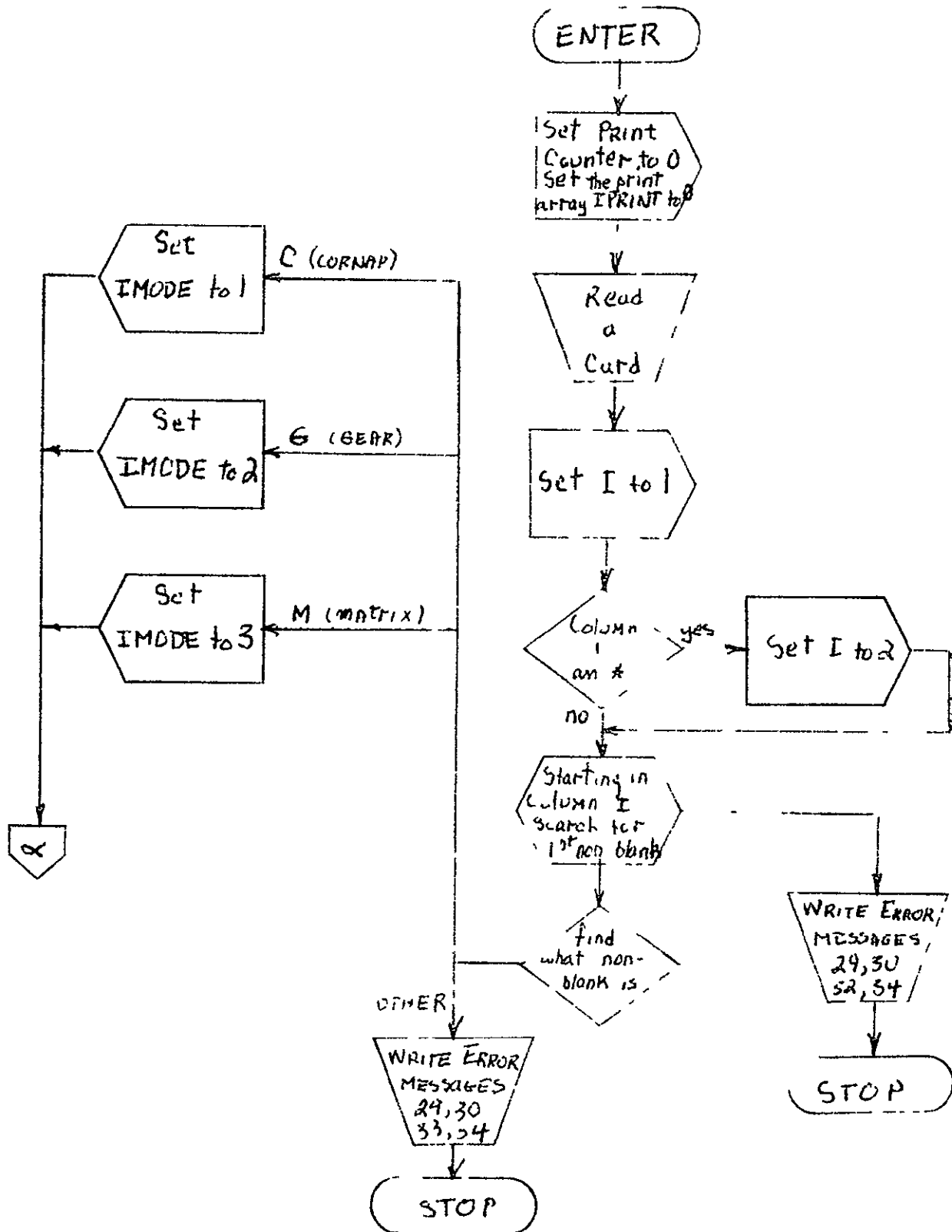


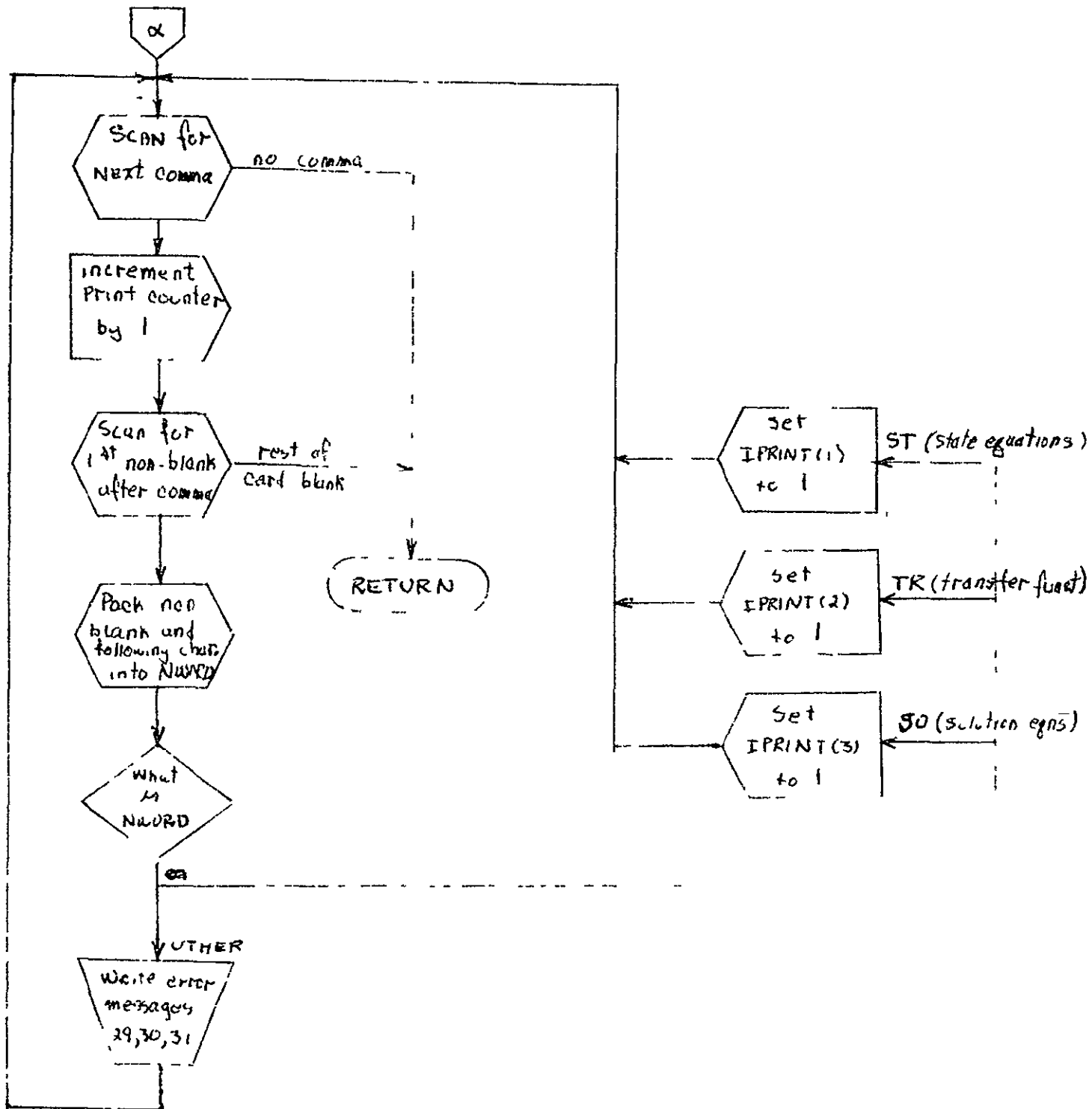


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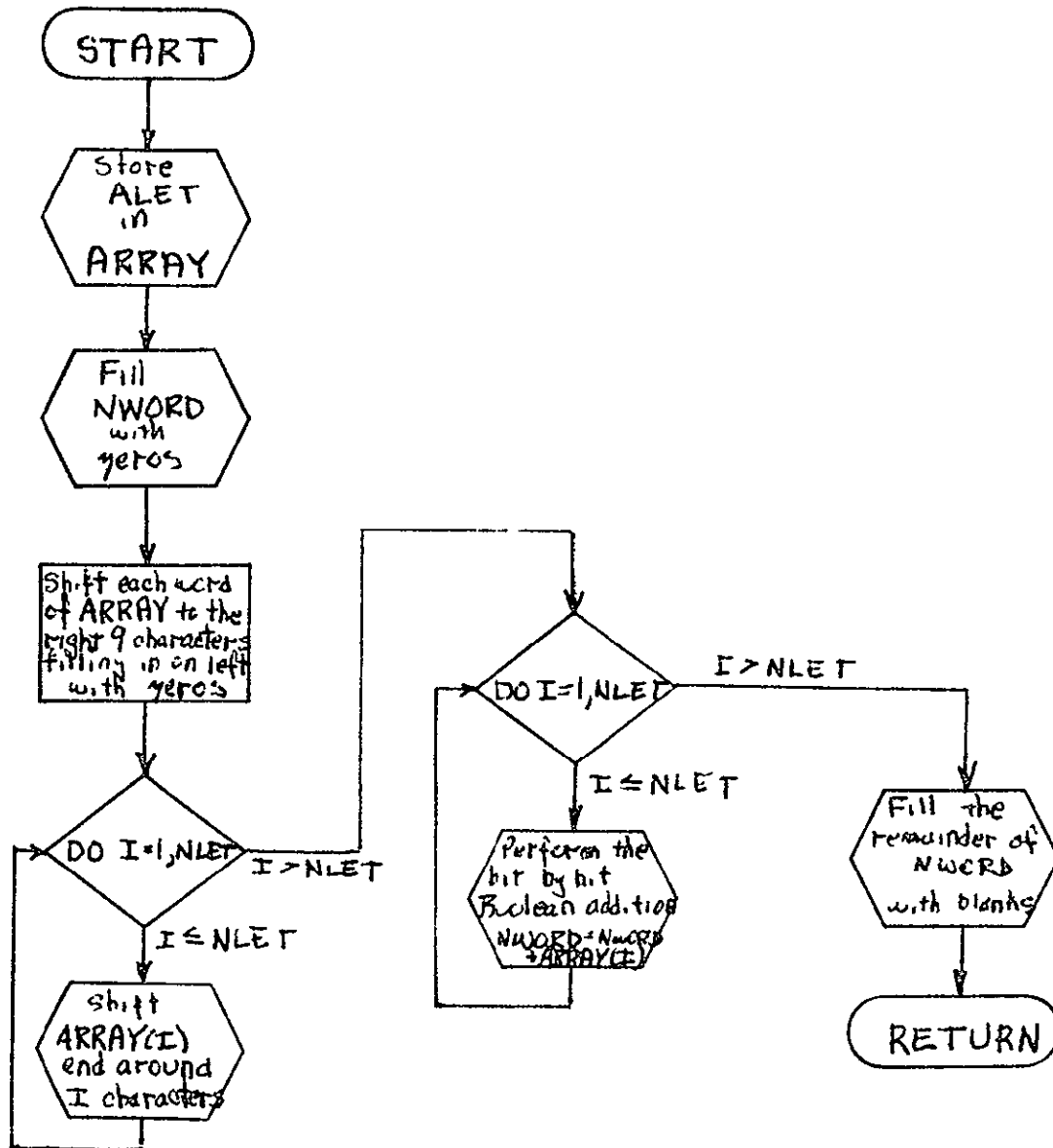


MODE



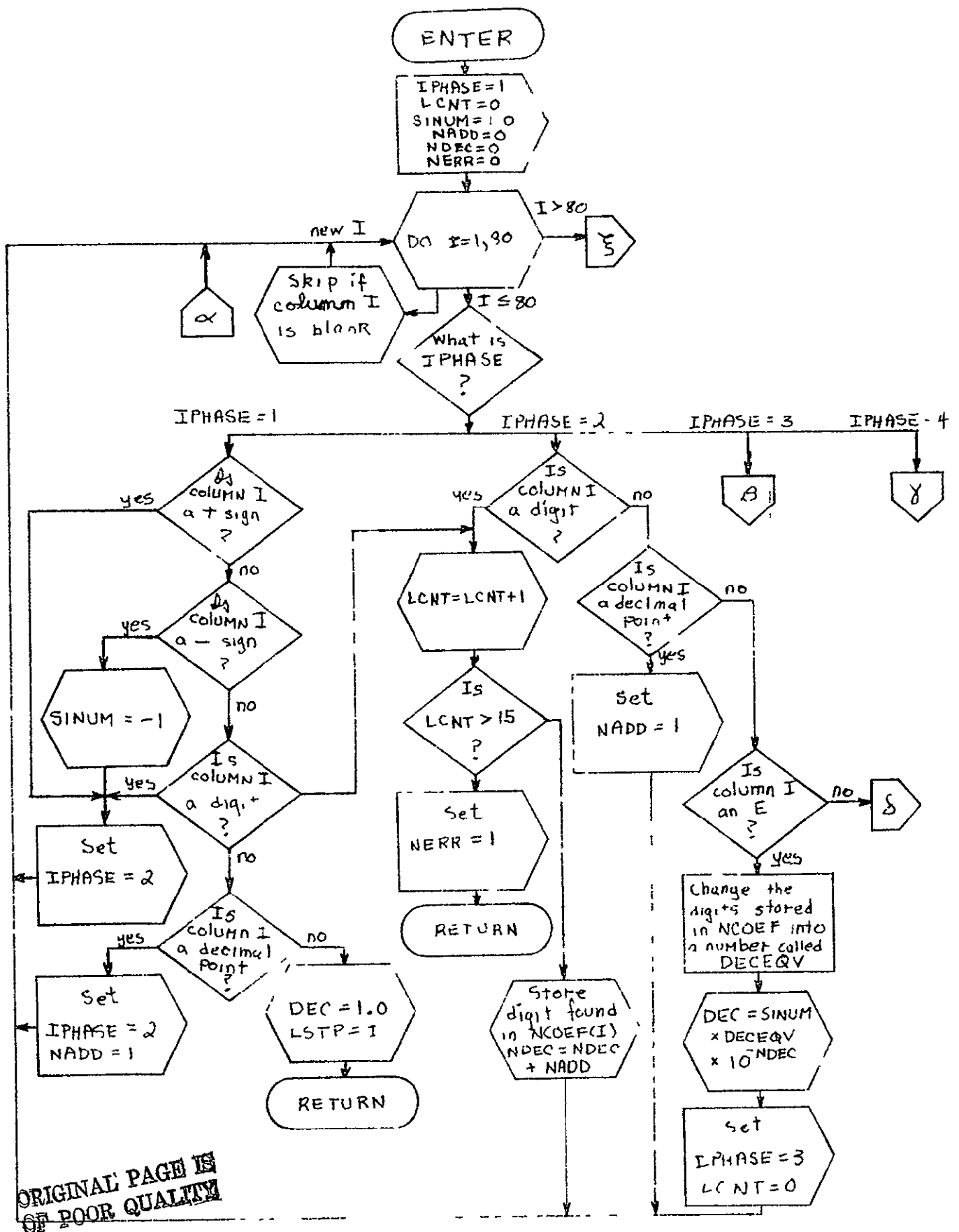


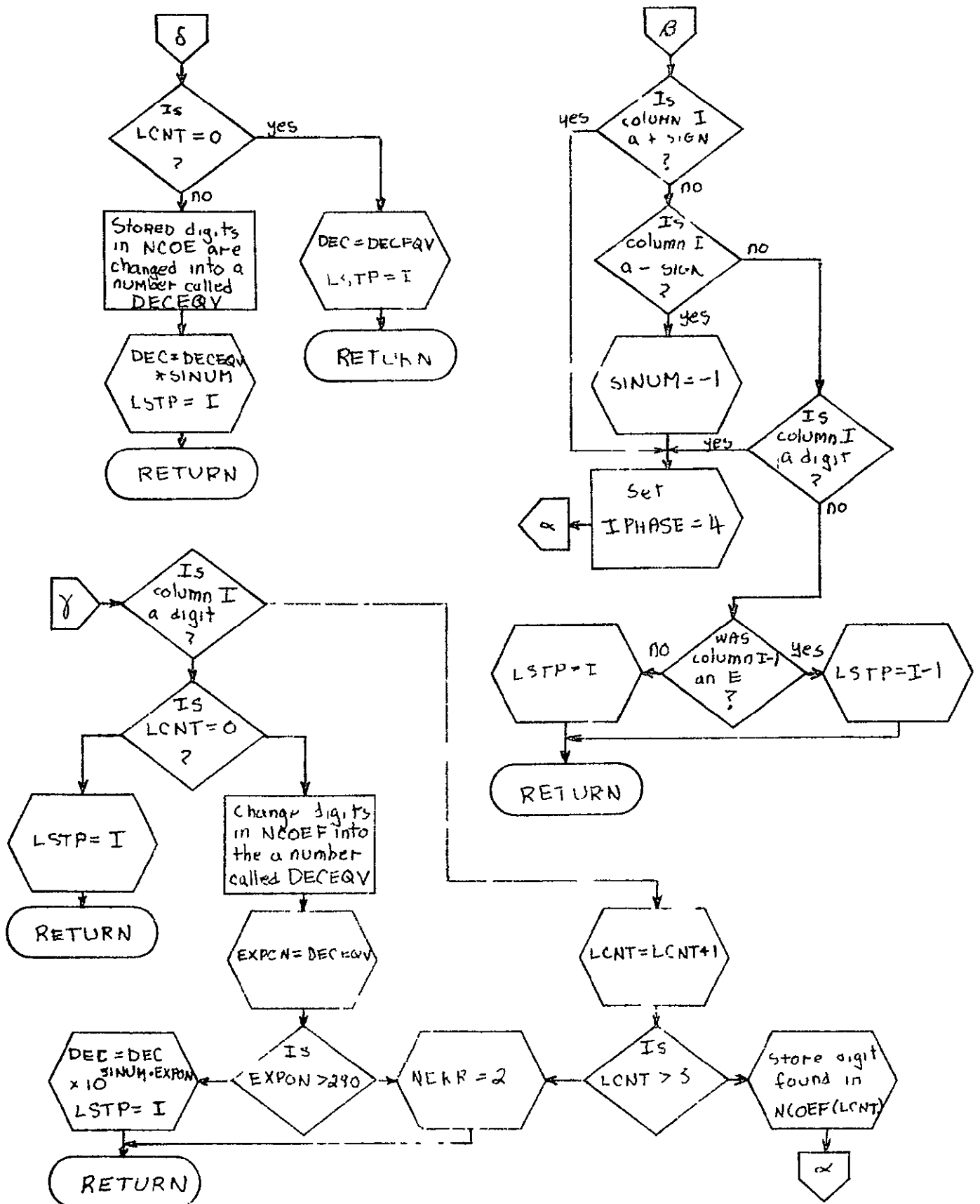
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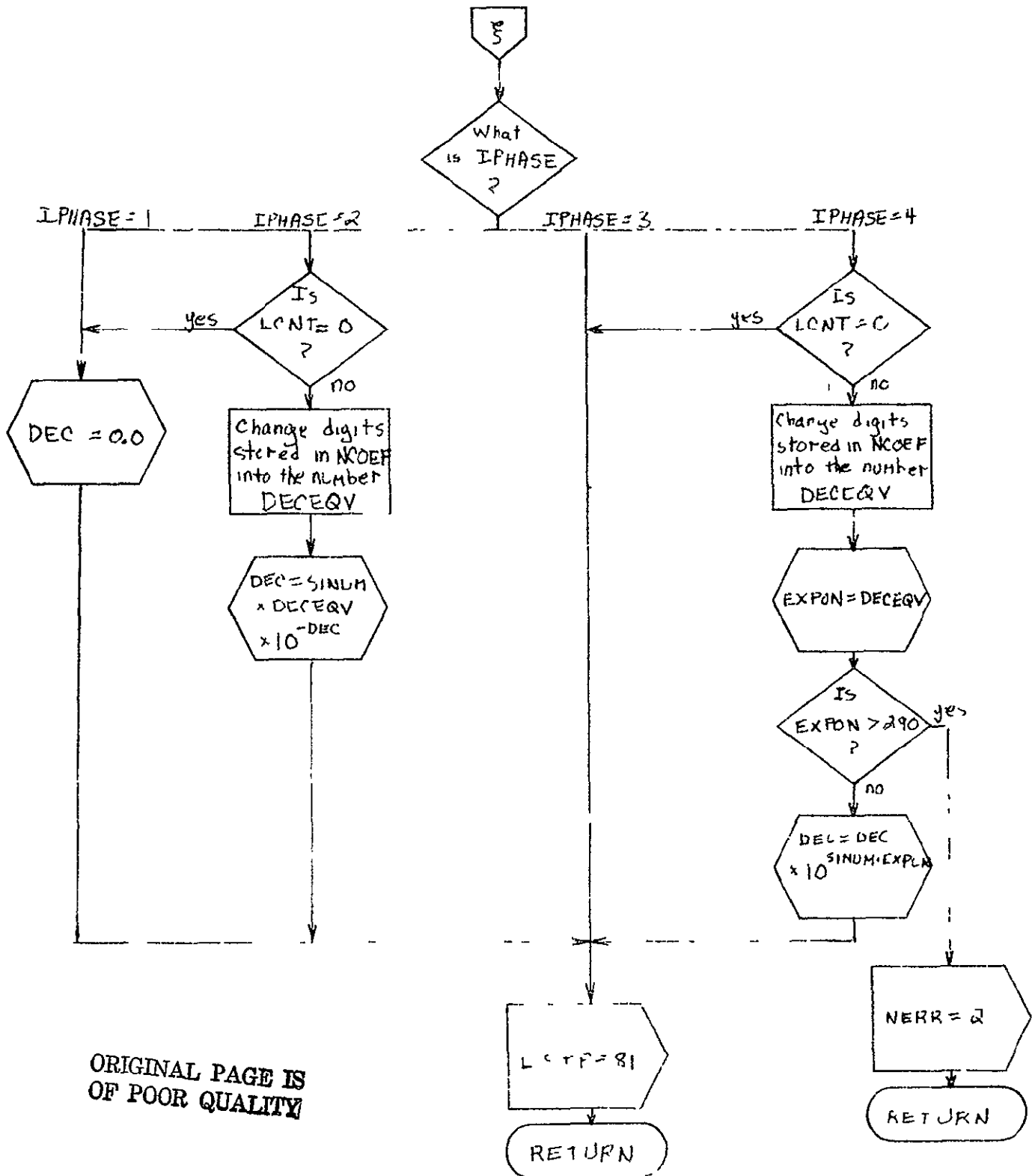


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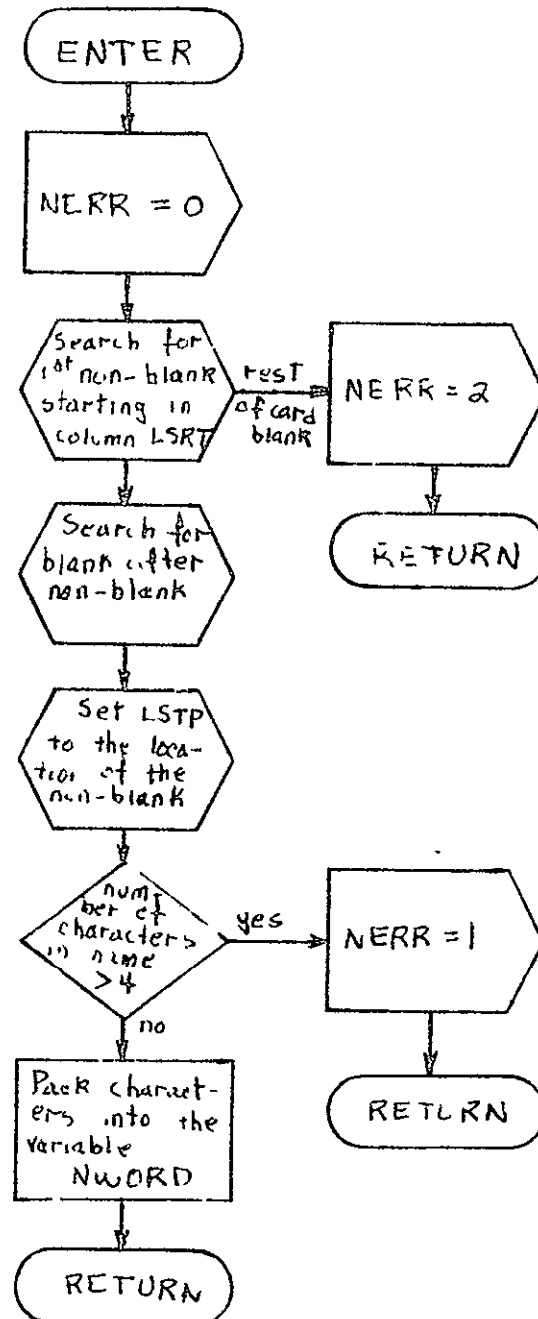
ATC DEC





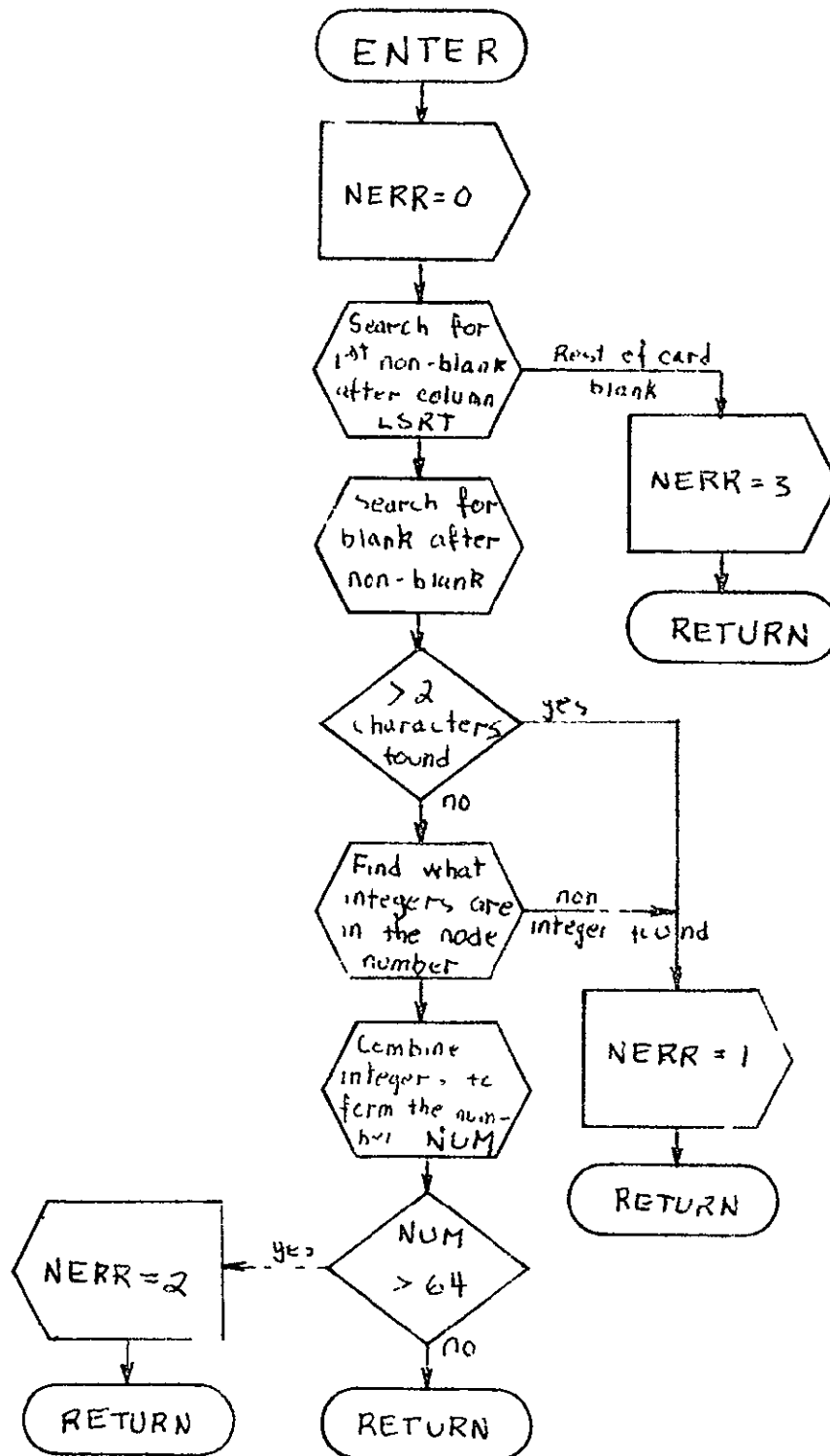


NAME

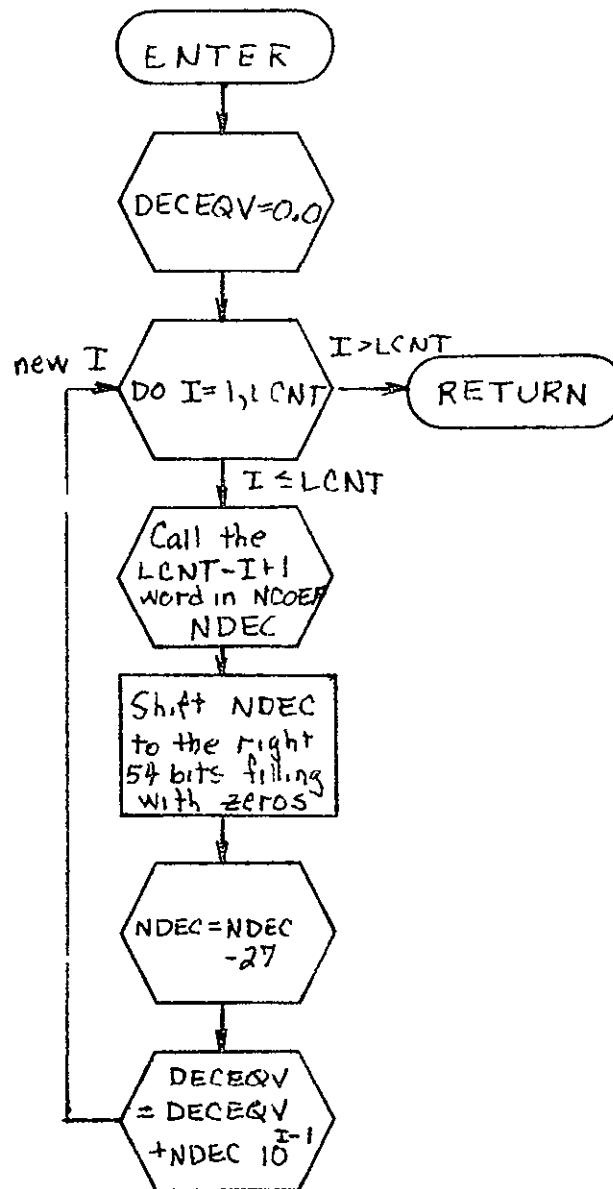


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NCDE

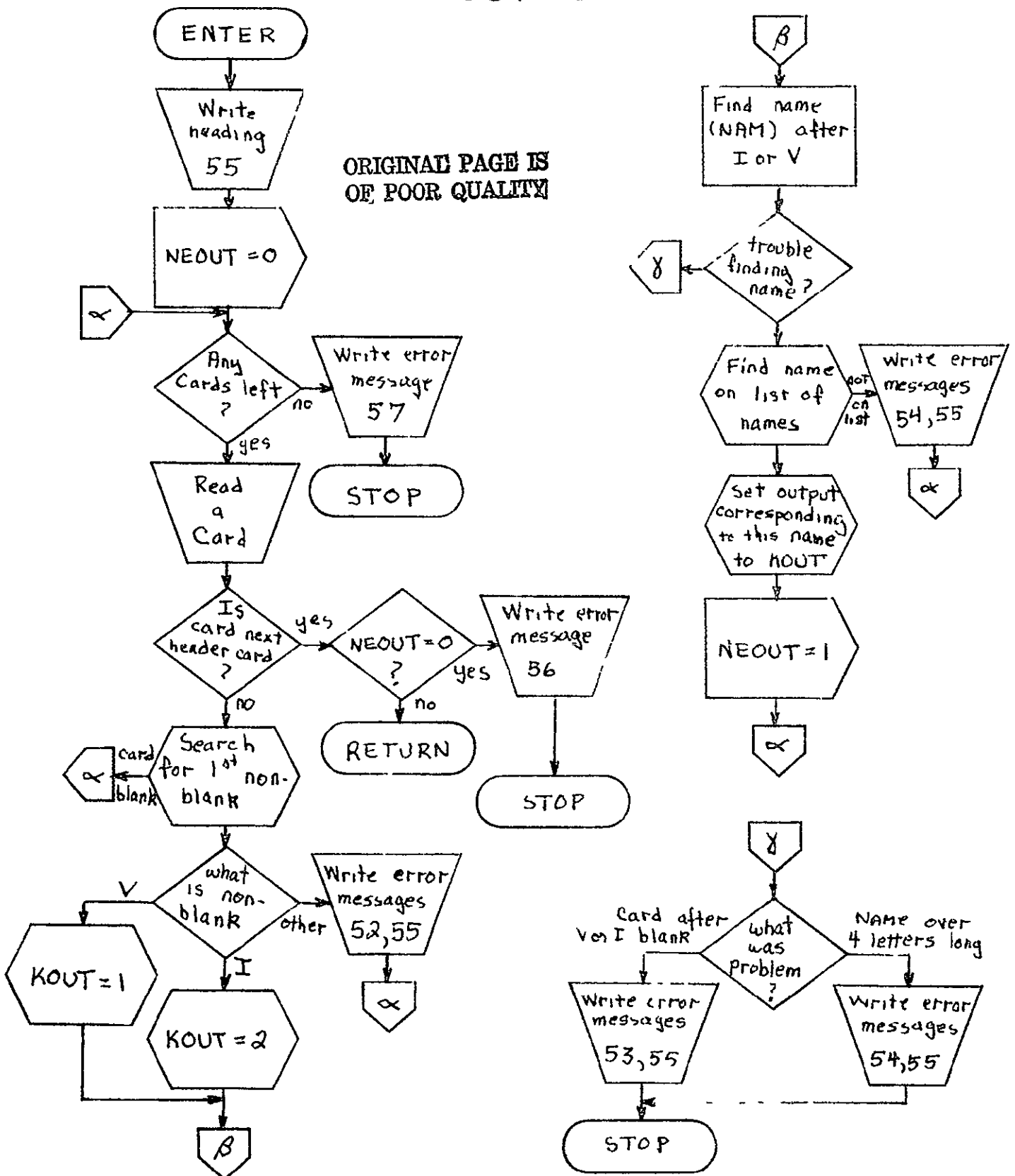


DECEQV

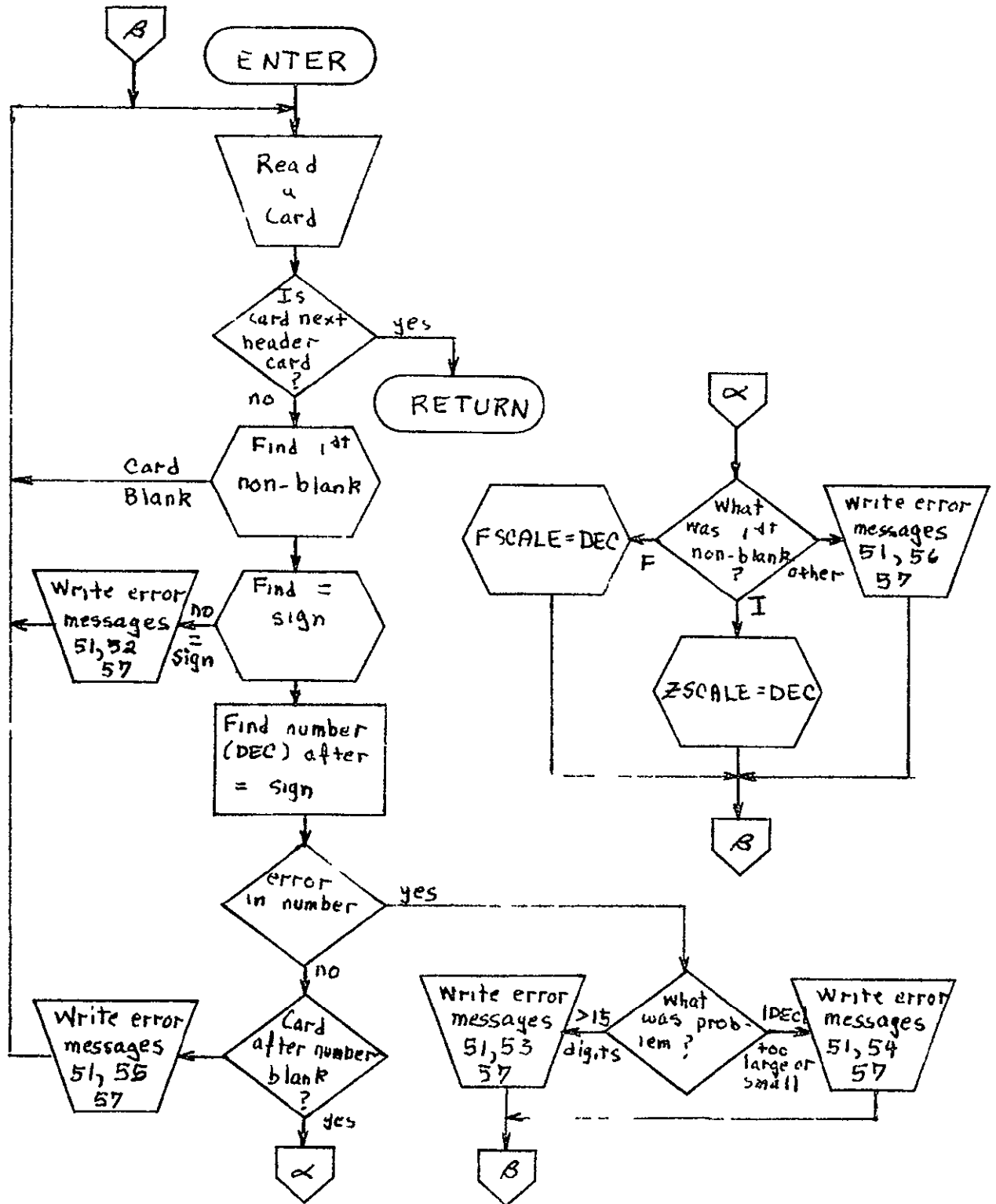


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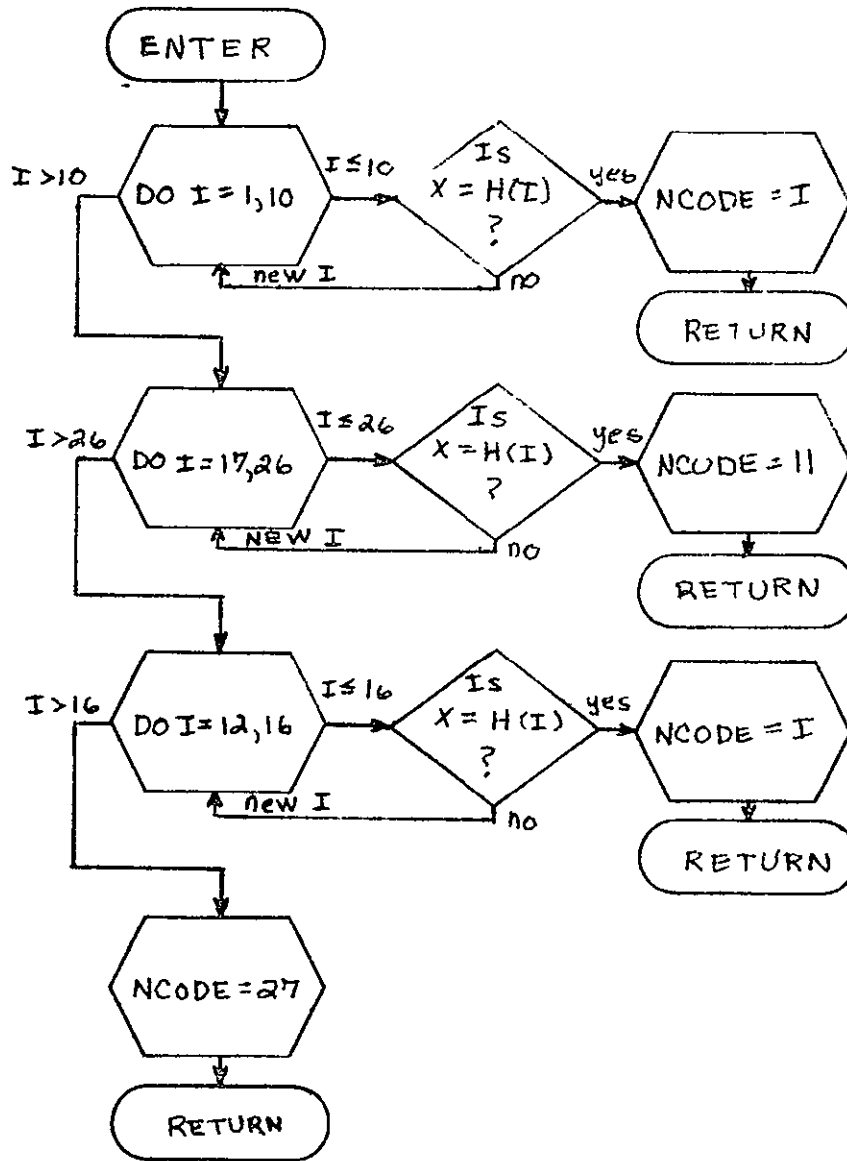
OUTPUT



Z F S C A L



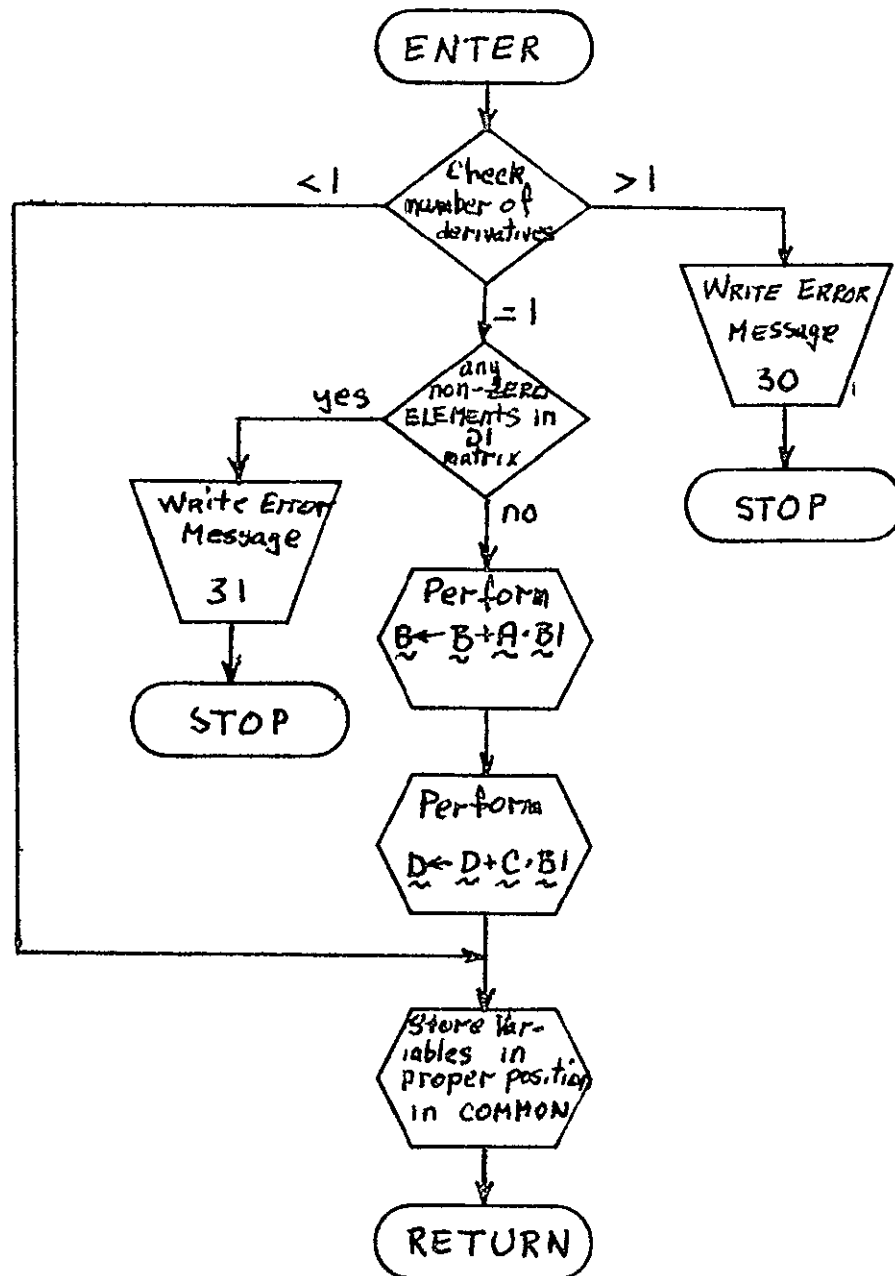
NCODE



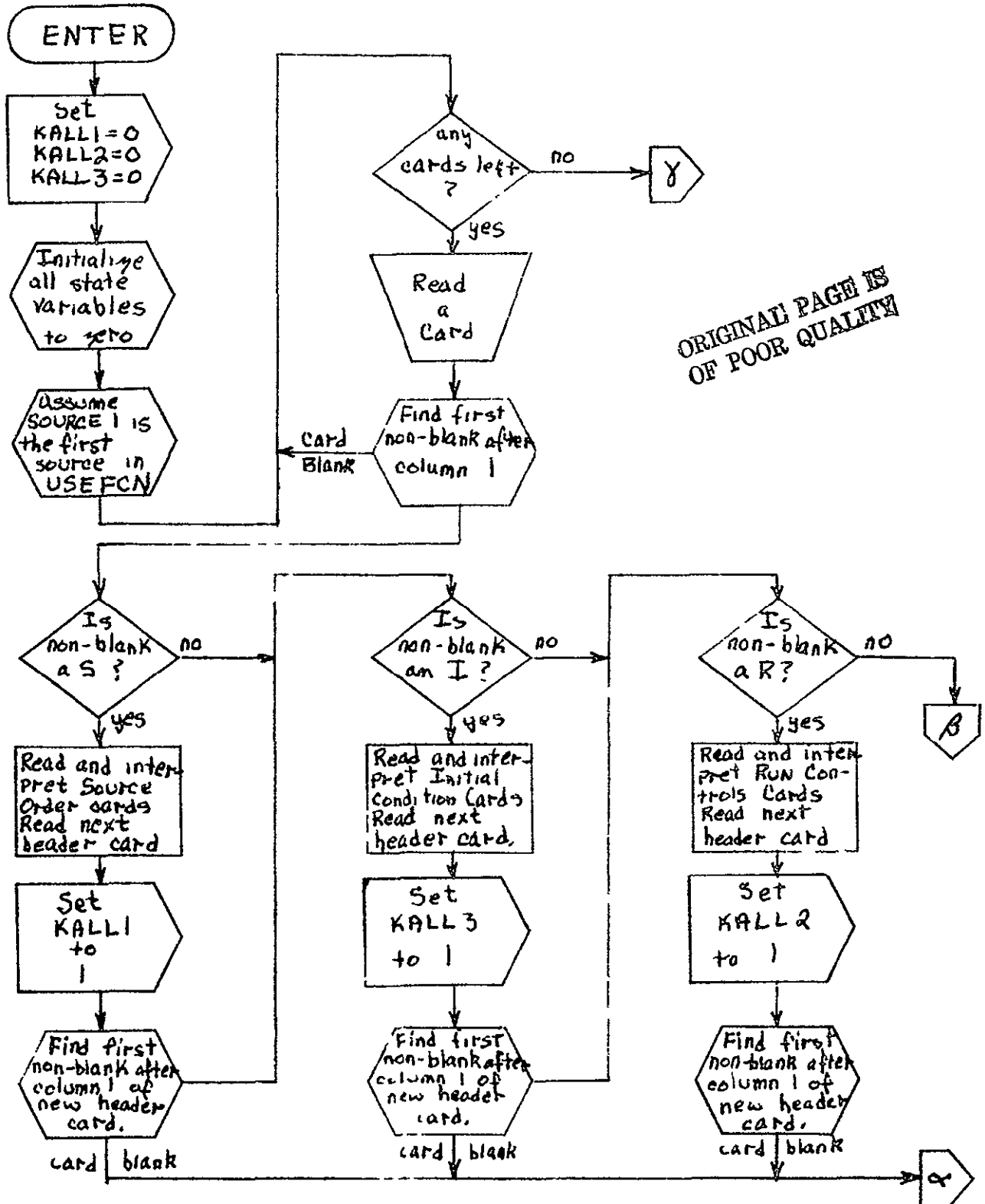
Here the array H contains

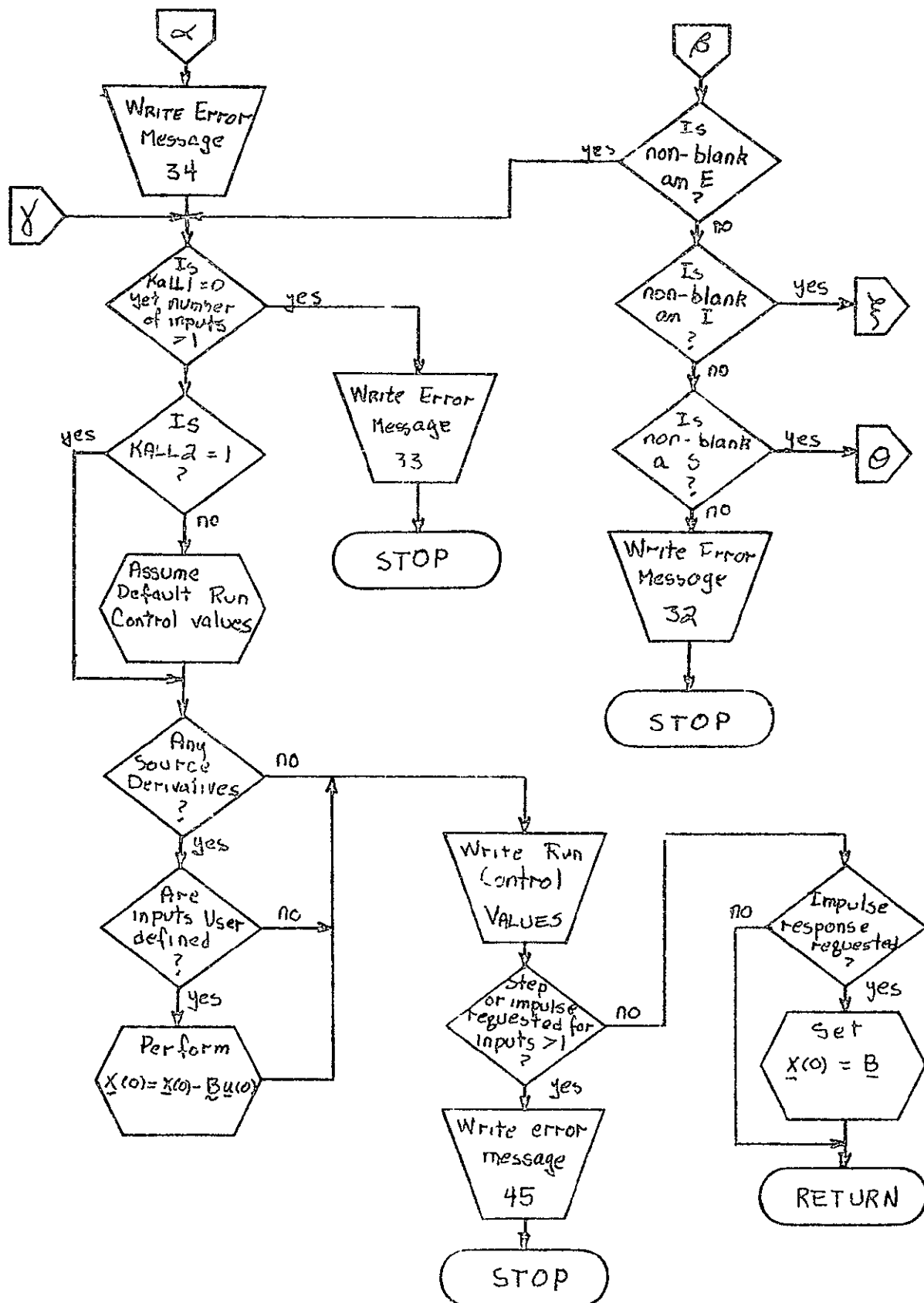
H(1)	1HV
H(2)	1HC
H(3)	1HR
H(4)	1HL
H(5)	1HI
H(6)	1H
H(7)	1H+
H(8)	1H-
H(9)	1HK
H(10)	1HM
H(11)	1H+
H(12)	1H(
H(13)	1H)
H(14)	1H
H(15)	1HE
H(16)	1H*
H(17)	1H1
H(18)	1H2
H(19)	1H3
H(20)	1H4
H(21)	1H5
H(22)	1H6
H(23)	1H7
H(24)	1H8
H(25)	1H9
H(26)	1H0

COMPAK

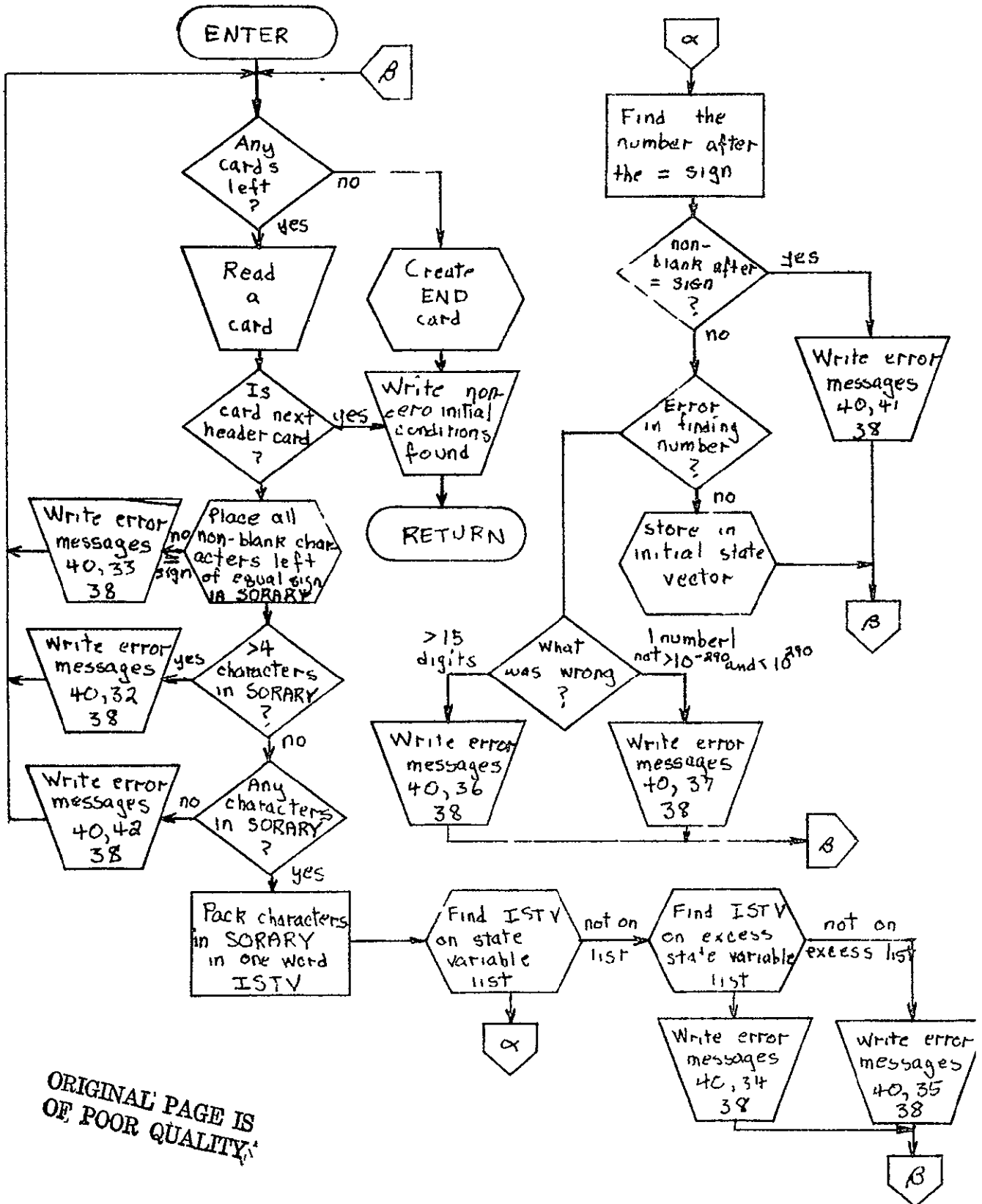


CROLGR



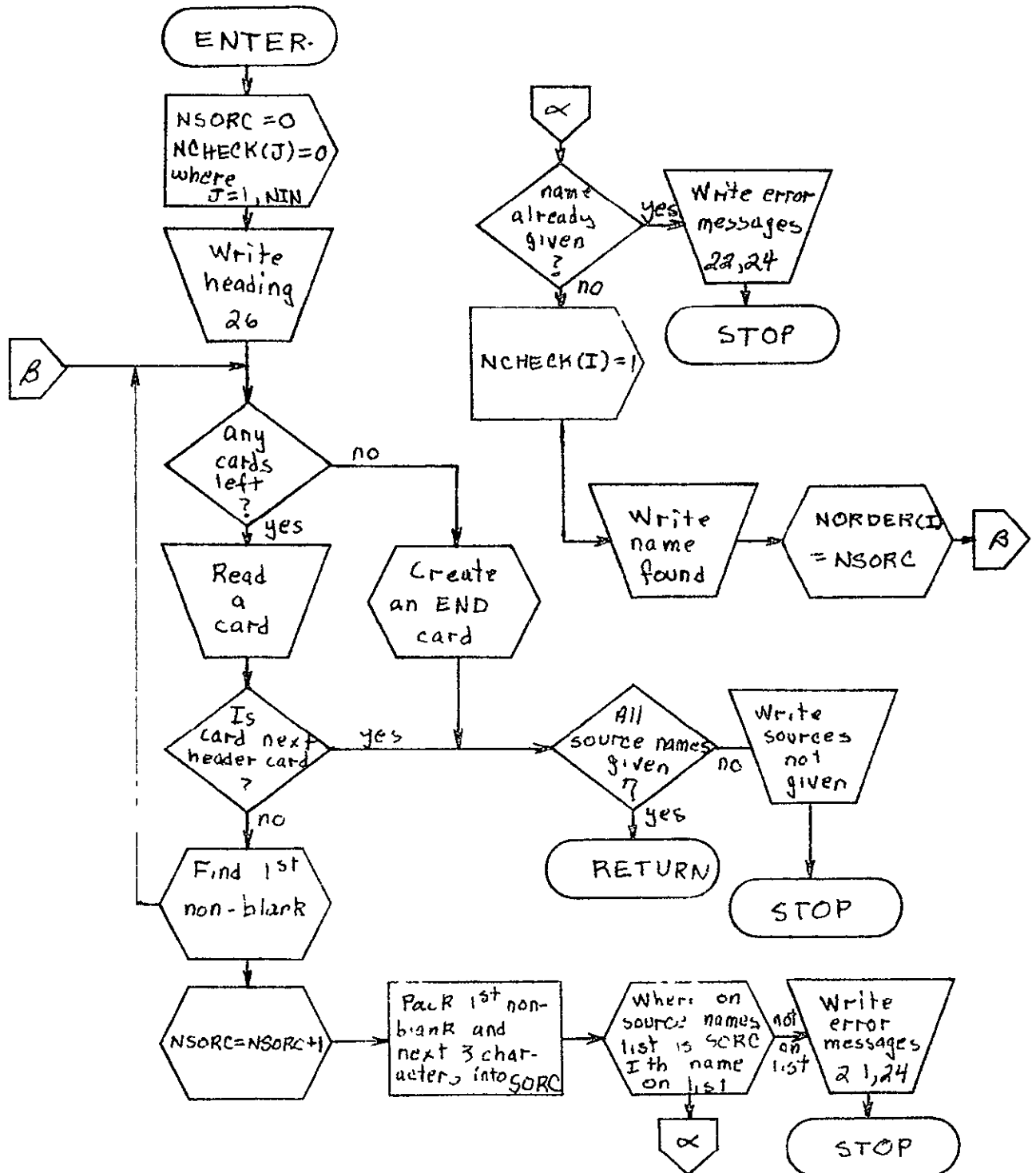


ICONS V

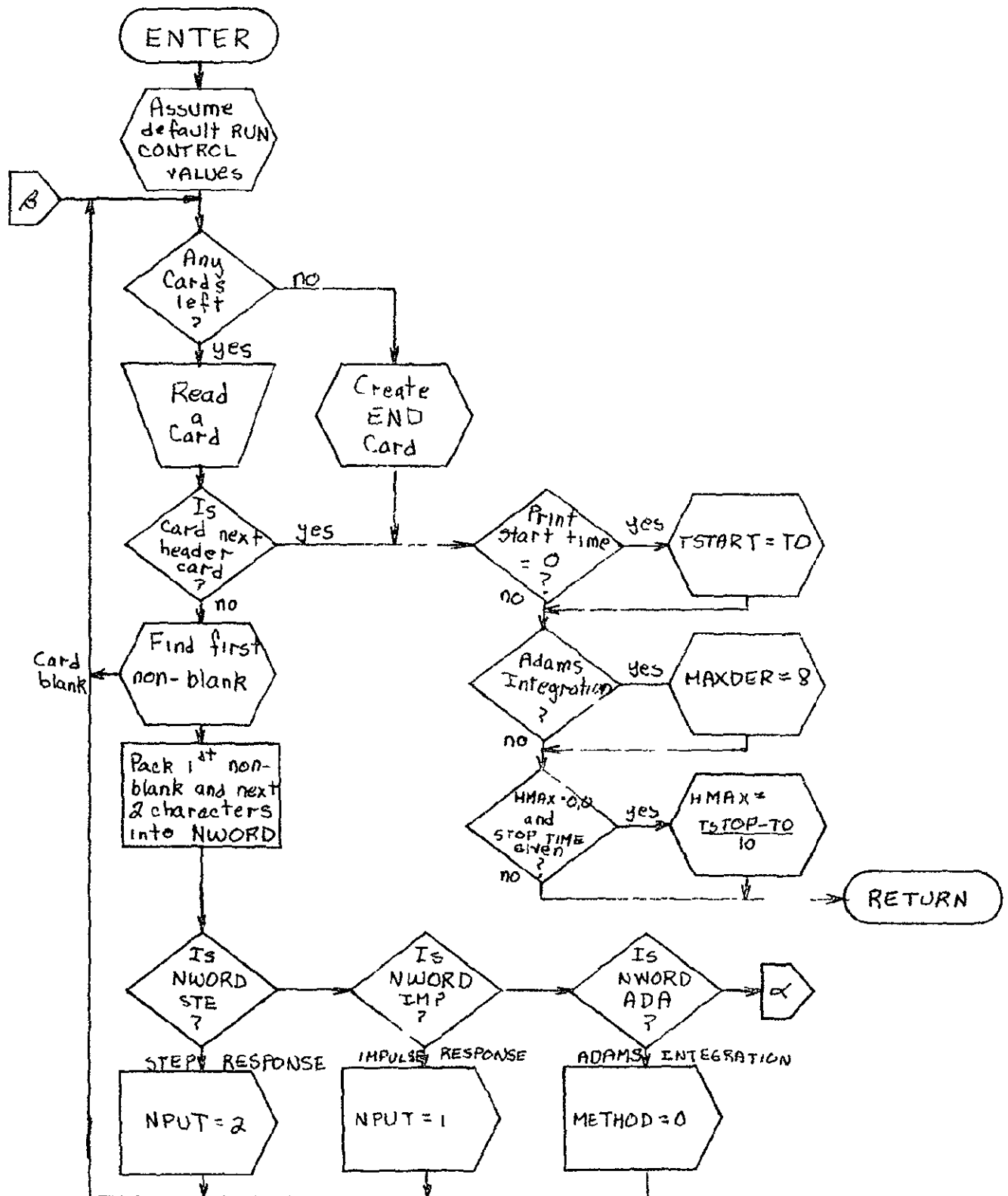


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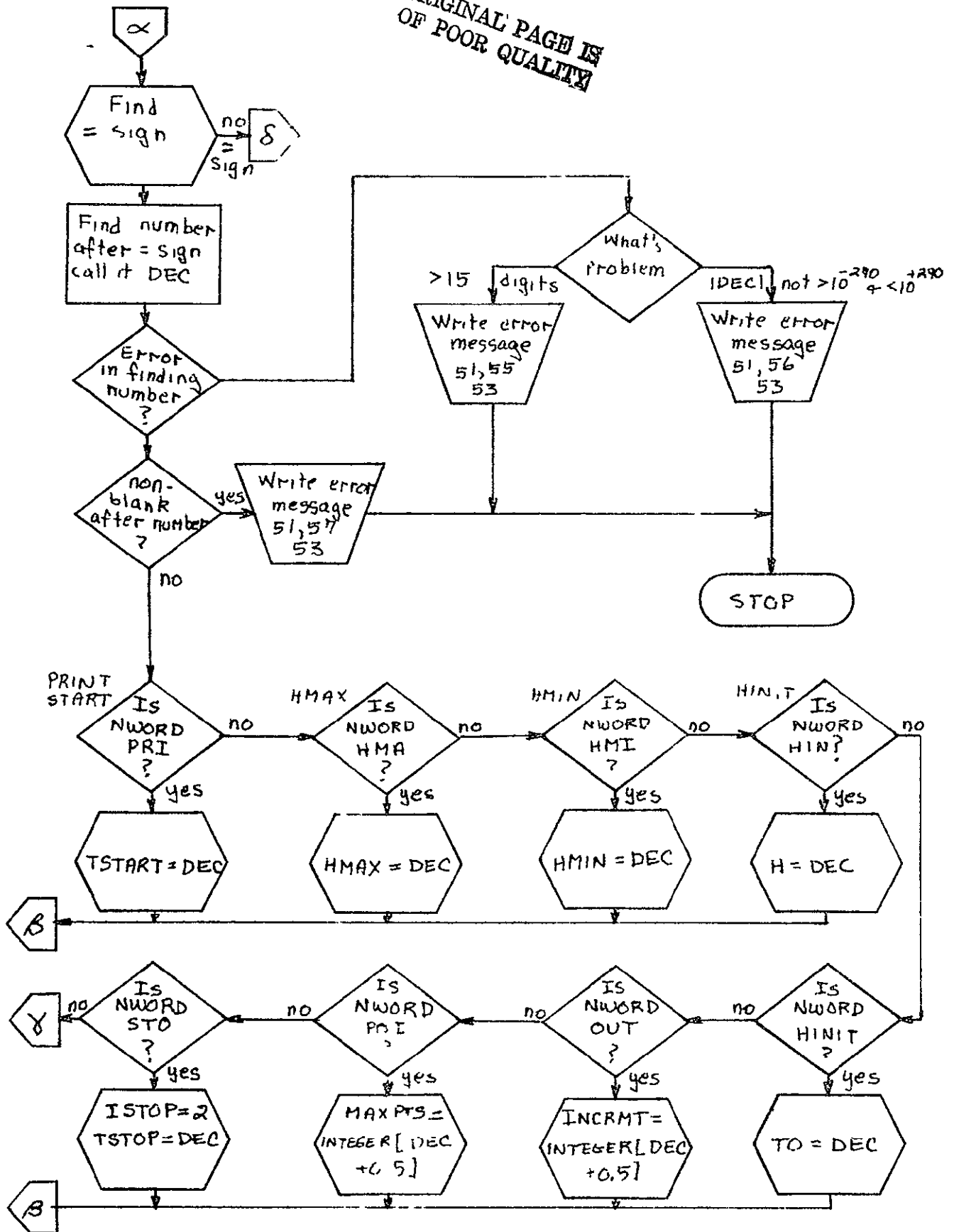
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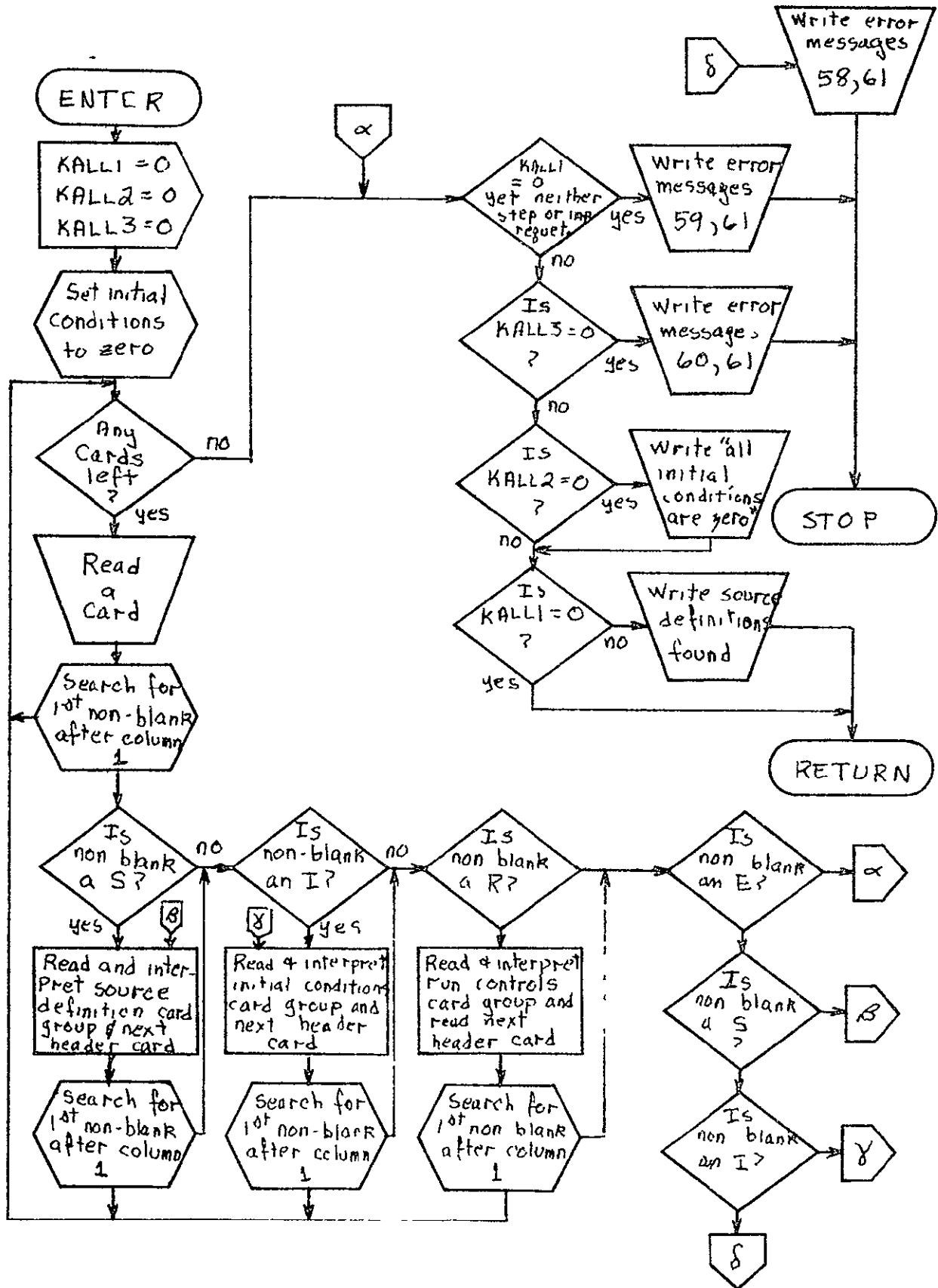
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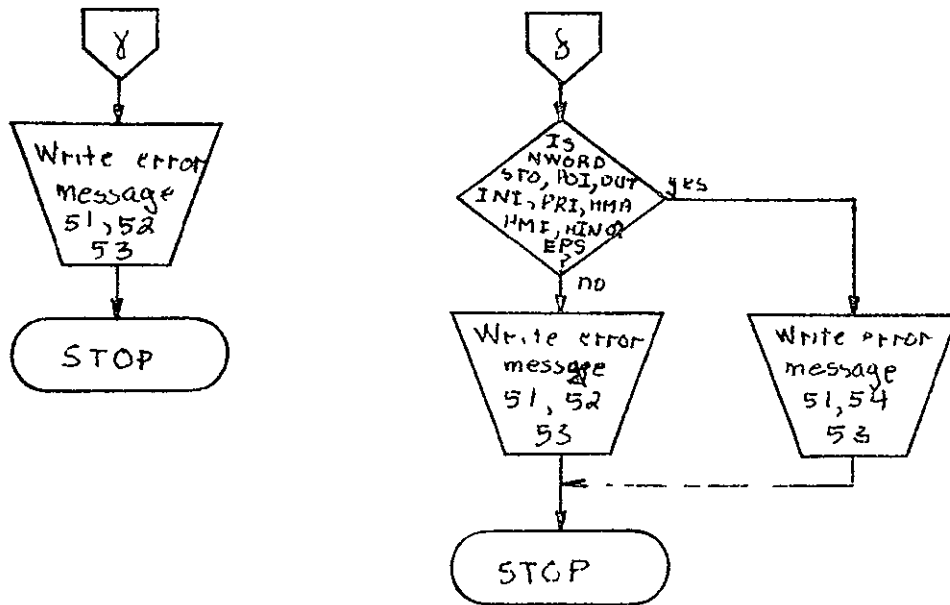


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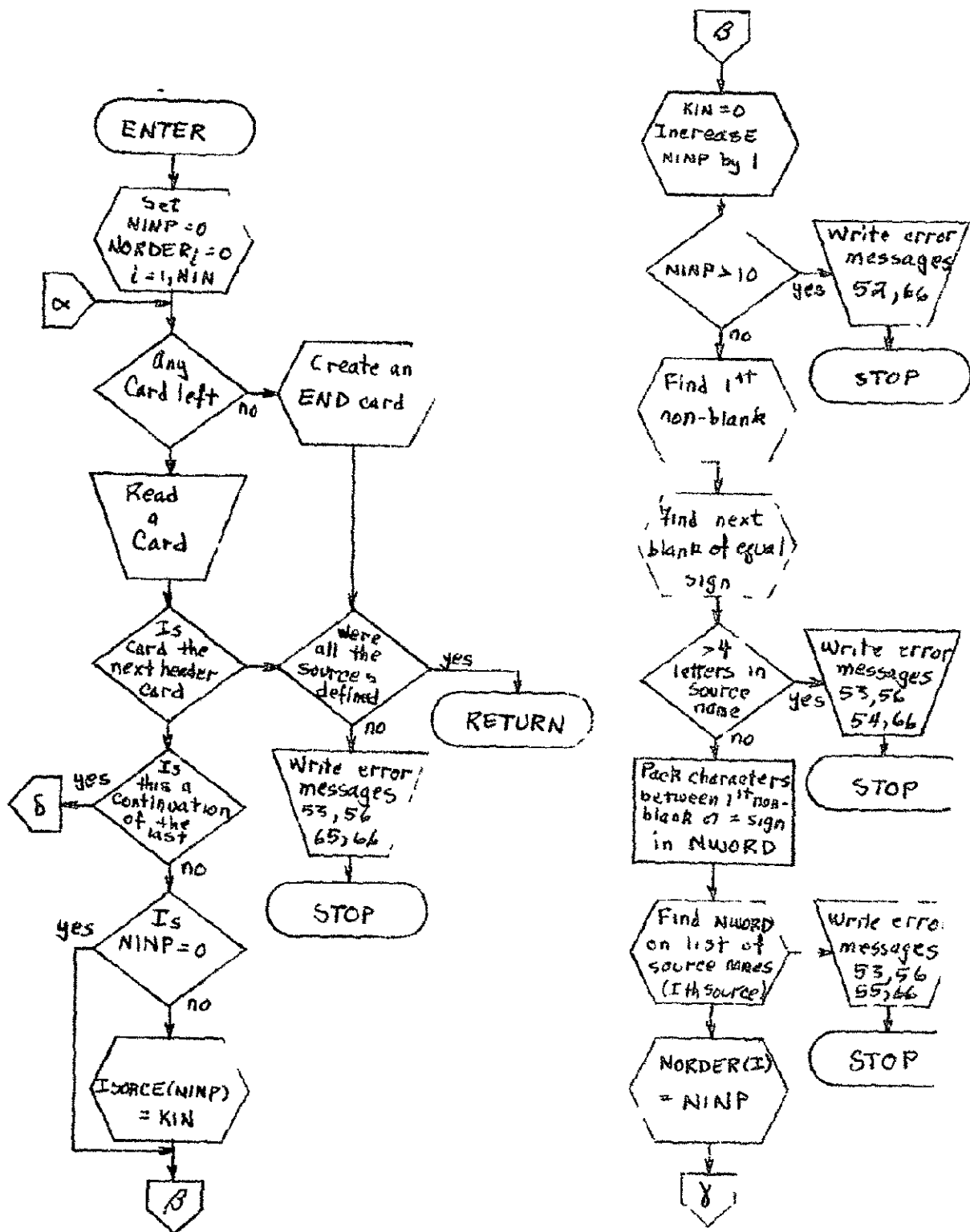
C R O L M X

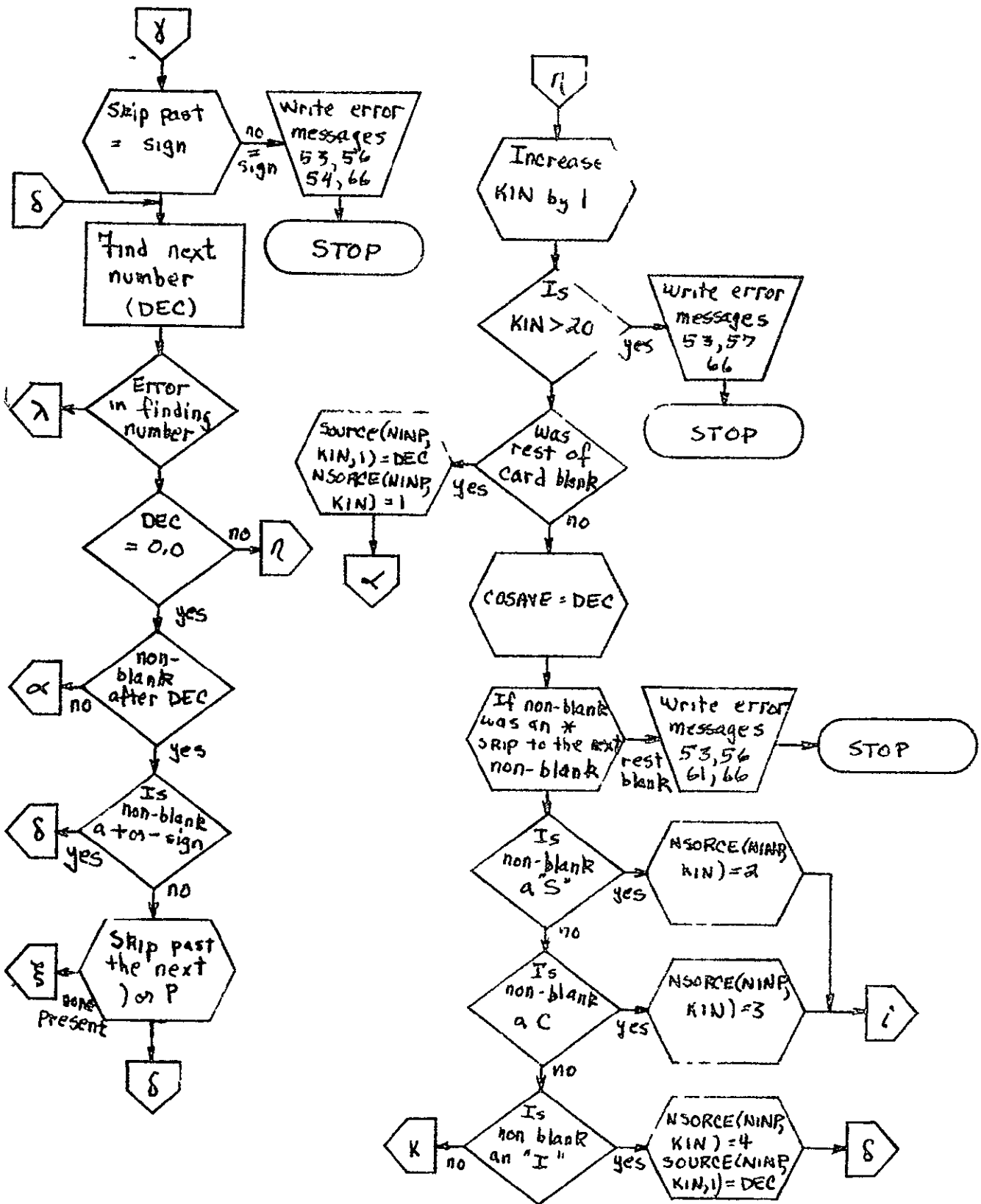




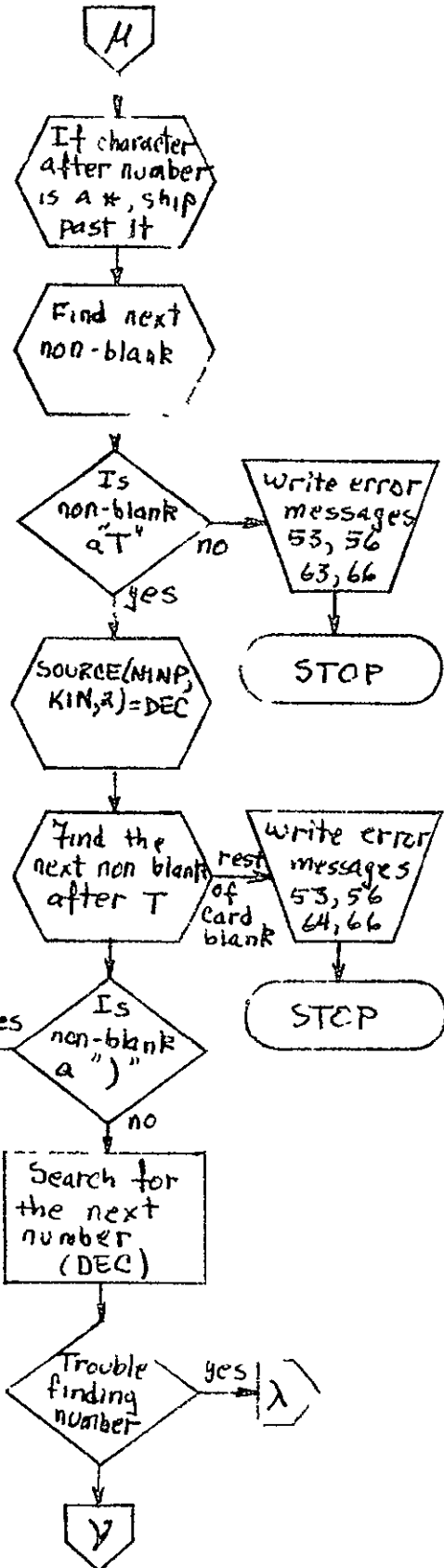
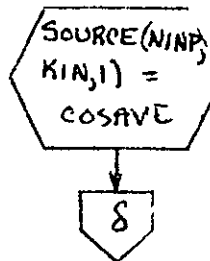
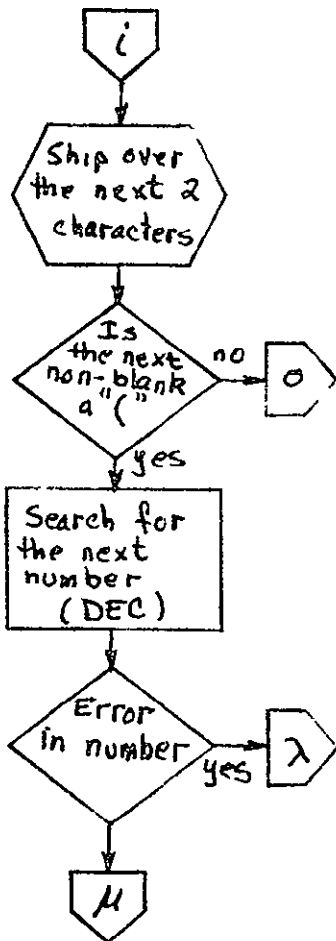
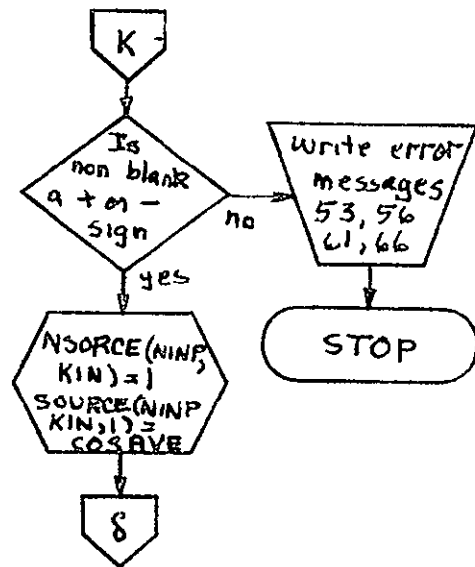
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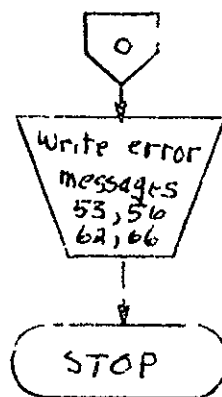
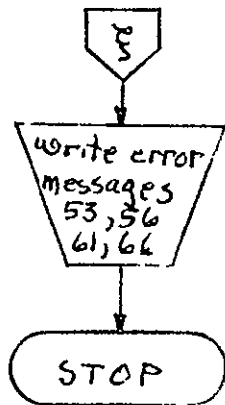
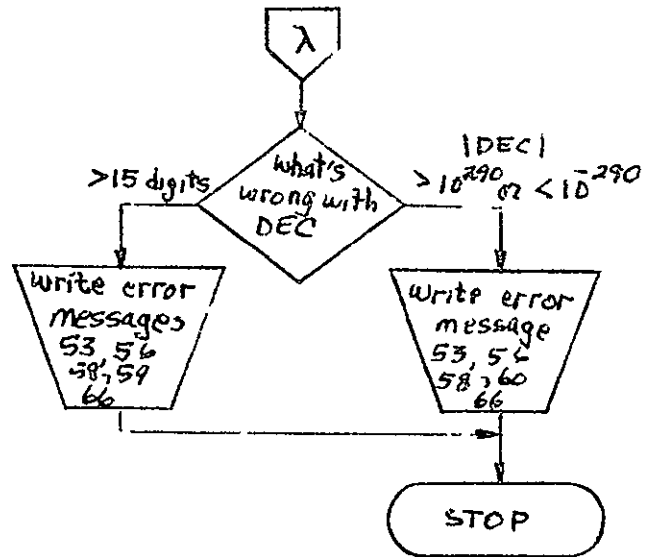
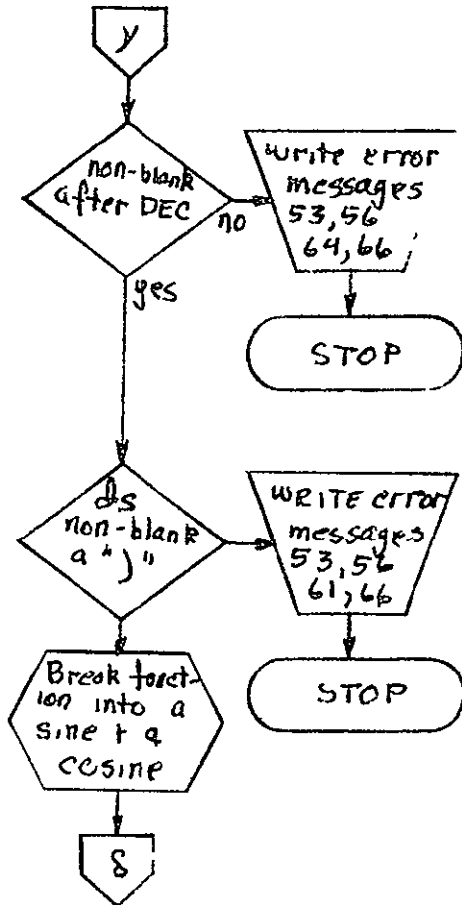
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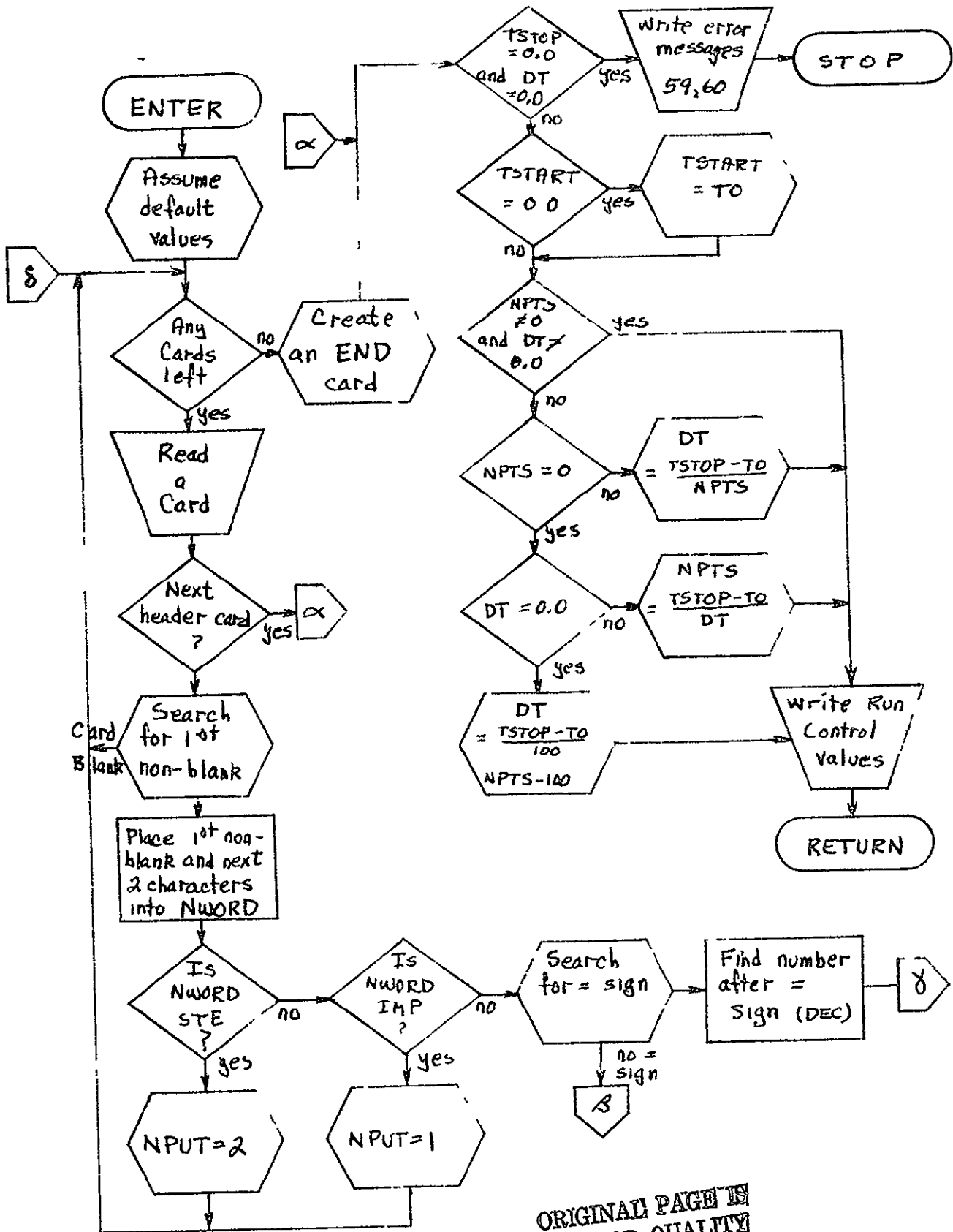


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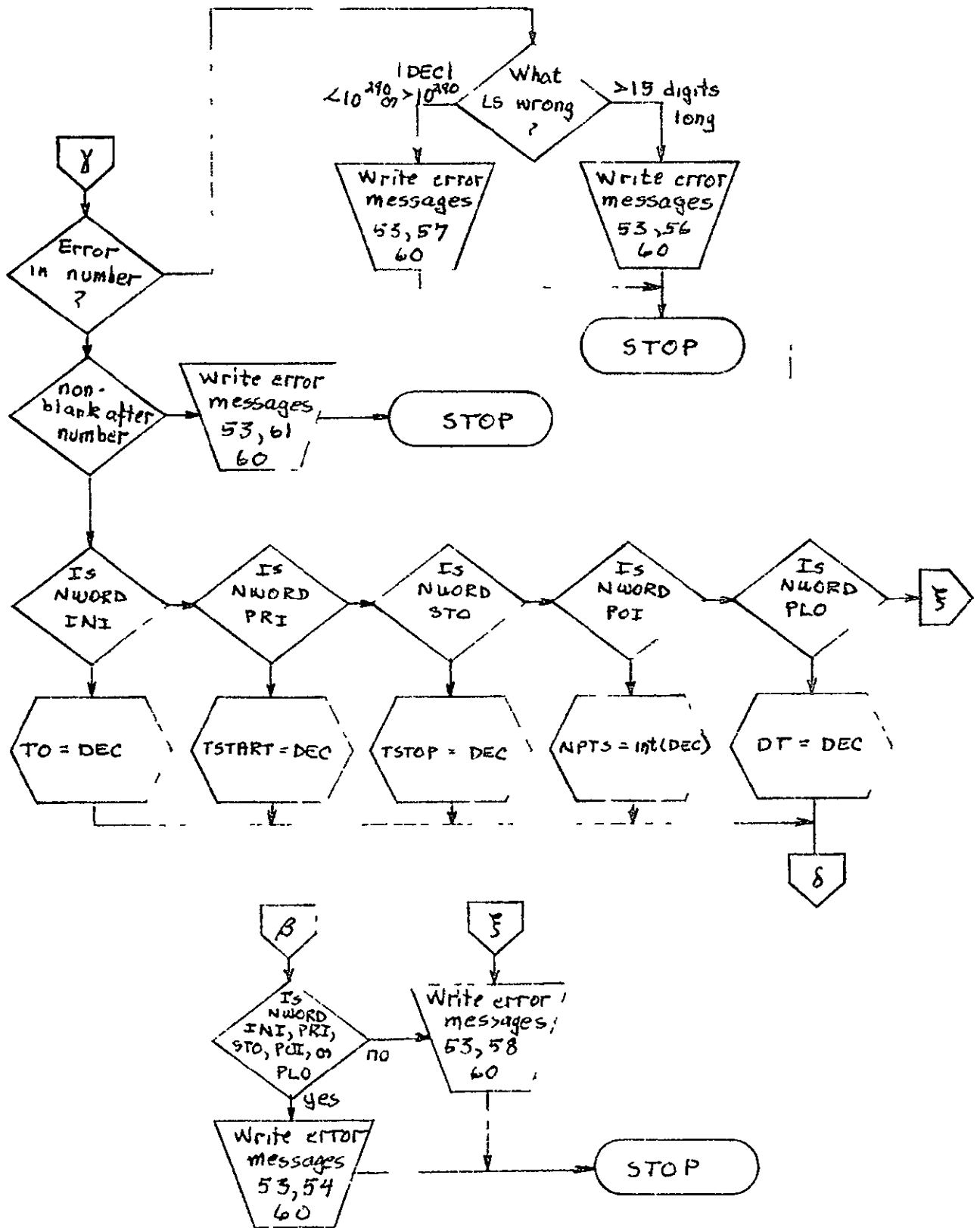




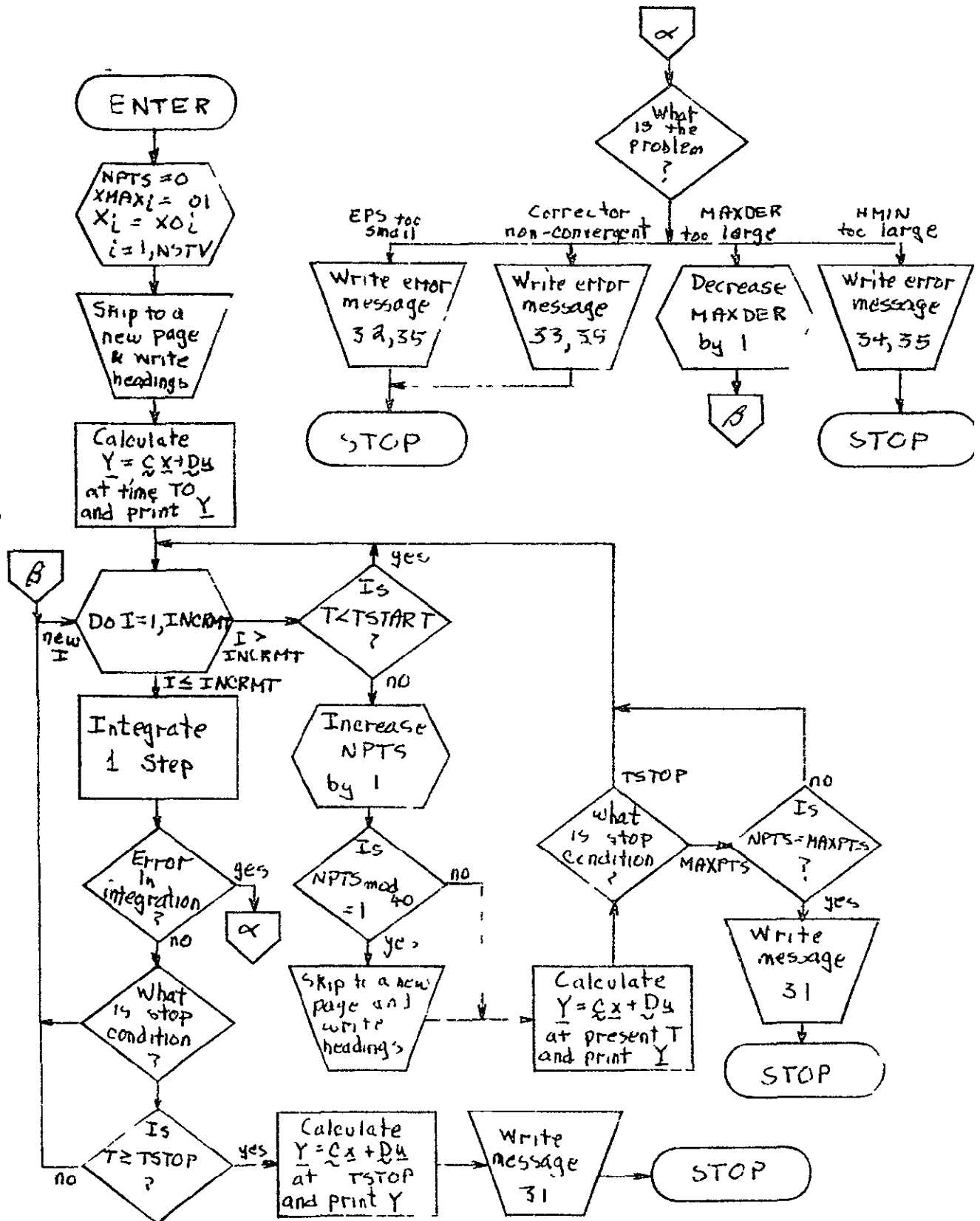
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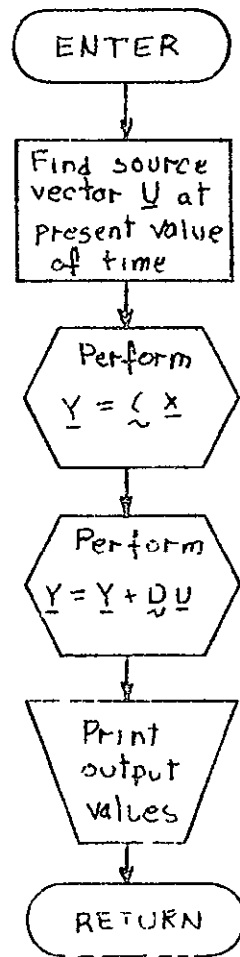
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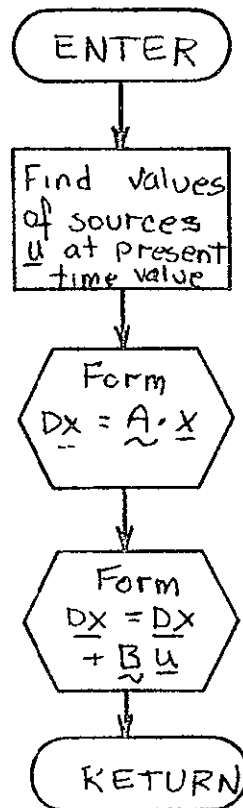
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DATPNT

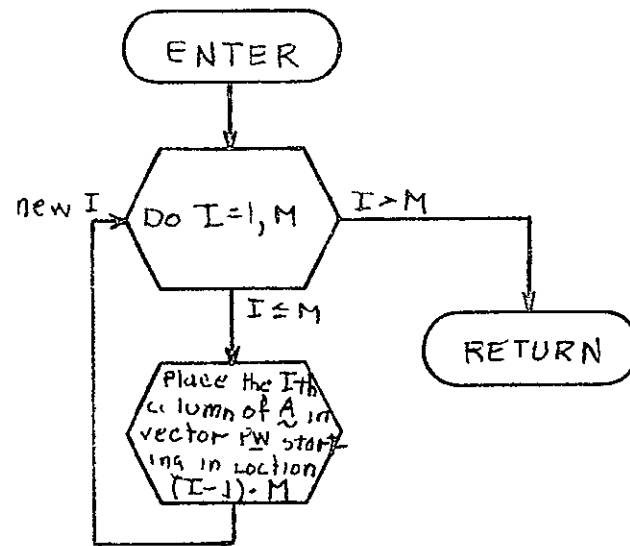


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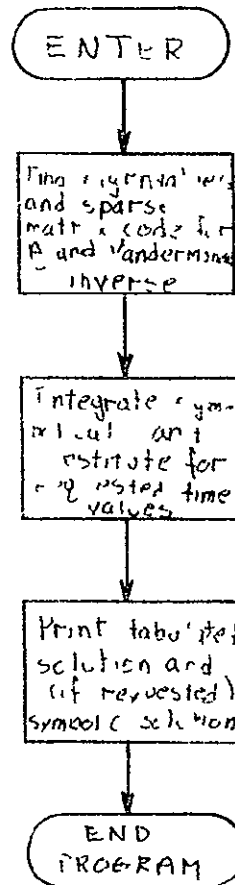


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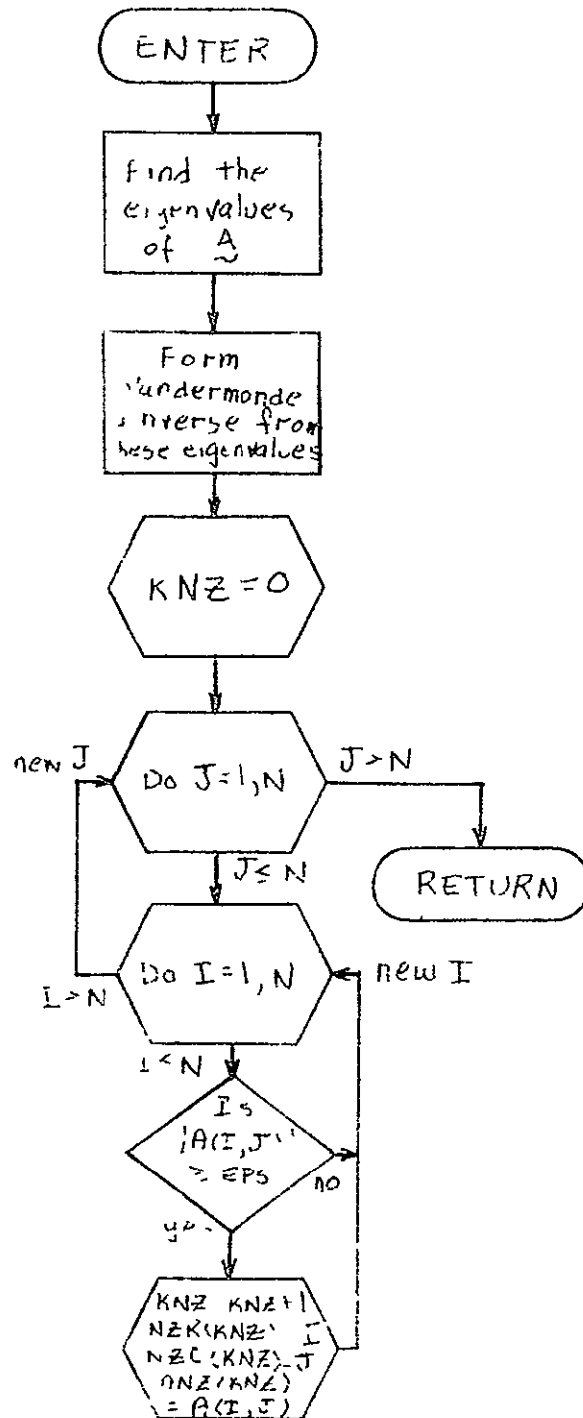


MATRIX

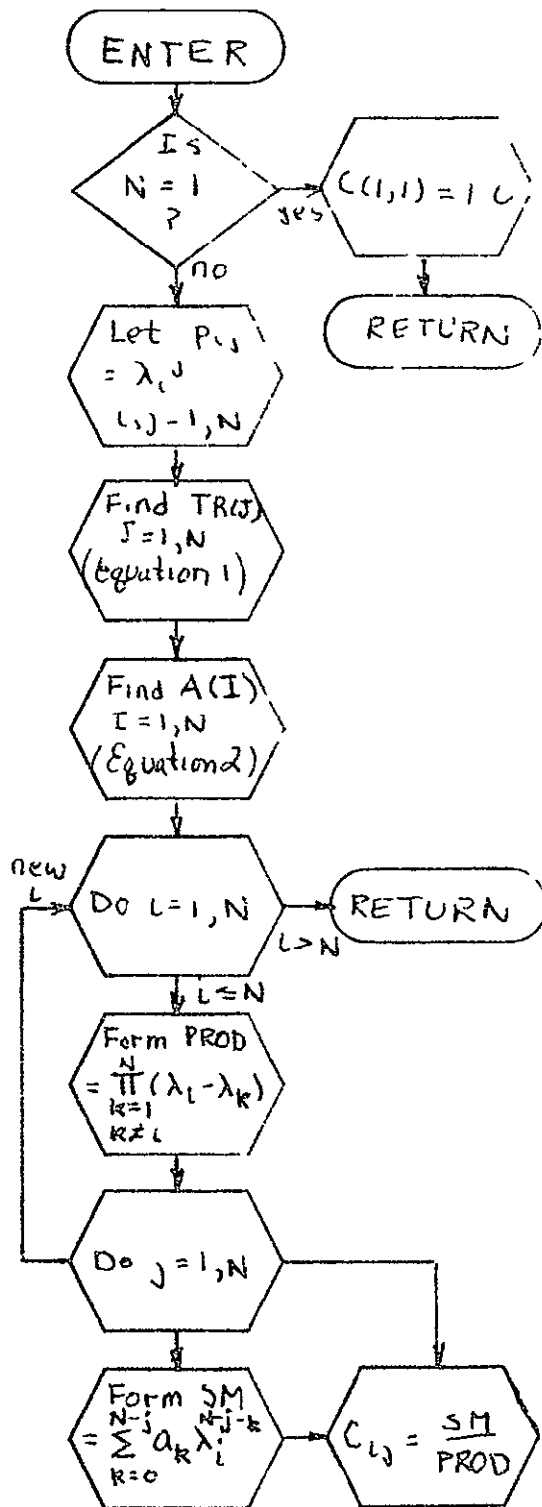


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SETSVE



VANINV



Equation 1

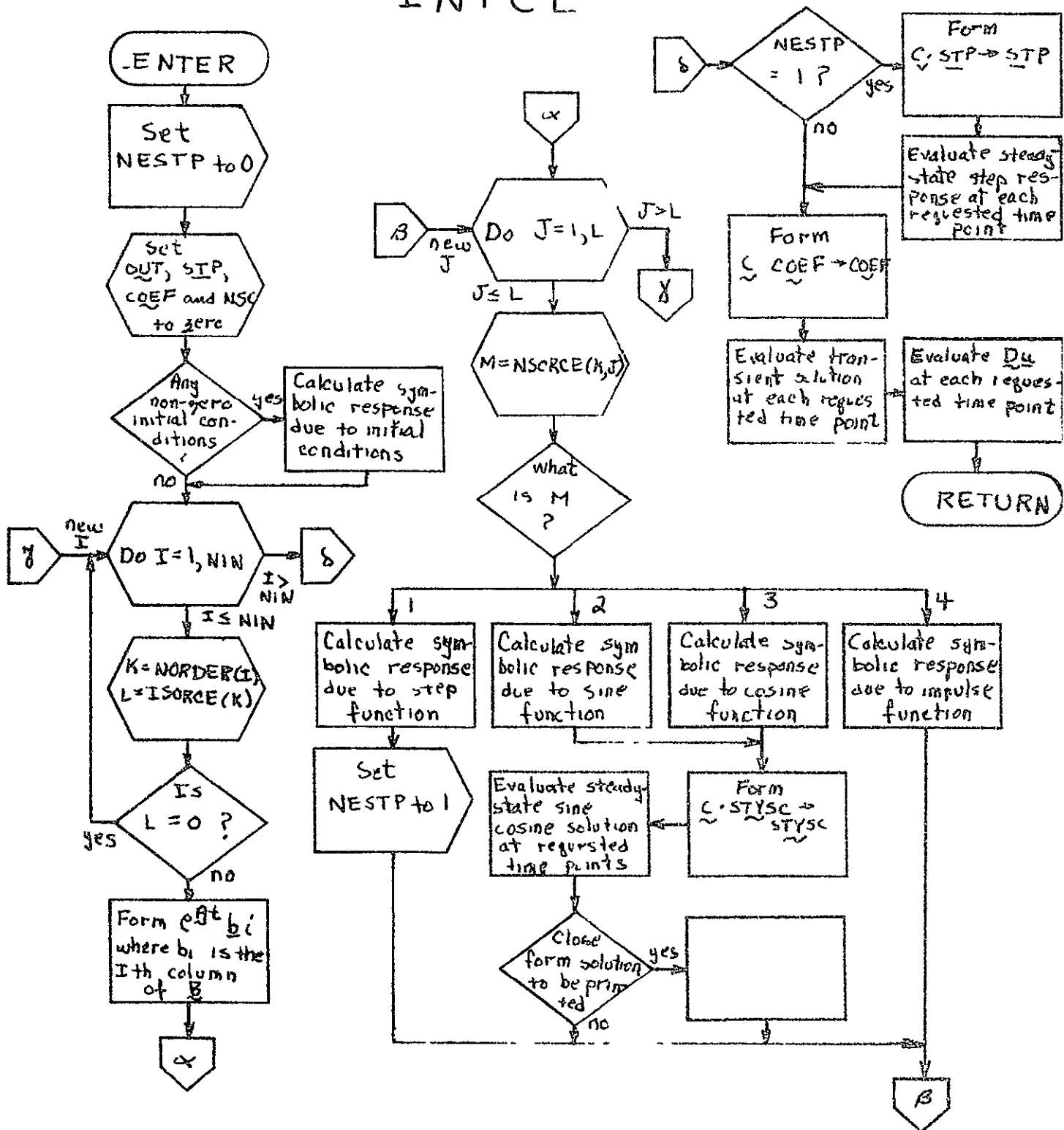
$$\begin{aligned} TR(J) &= TR A^J \\ &= \sum_{I=1}^N \lambda_I^J \\ &= \sum_{I=1}^N P_{J,I} \end{aligned}$$

Equation 2

$$|A - \lambda I| = \lambda^n + a_1 \lambda^{n-1} + \dots + a_n$$

$$A(I) = a_i = -\frac{1}{i} \sum_{m=1}^i a_m TR(1m)$$

INTCL



SVECOS

Equation 1

$$B_i = \text{ROOT}_i^2 + \text{OMEGA}^2$$

$$C_i = \text{AMPL} \cdot e^{-\text{ROOT}_i \cdot \text{TO}} [-\text{ROOT}_i \cdot \cos(\text{OMEGA} \cdot \text{TO}) + \text{OMEGA} \sin(\text{OMEGA} \cdot \text{TO})] / B_i$$

$$D_i = -\text{AMPL} \cdot \text{ROOT}_i / B_i$$

$$E_i = \text{AMPL} \cdot \text{OMEGA} / B_i$$

$$\text{STYCOS}_{j,1} = \text{STYCOS}_{j,1} + P_{j,1} \cdot E_i$$

$$\text{STYCOS}_{j,2} = \text{STYCOS}_{j,2} + P_{j,2} \cdot D_i$$

$$\text{CFCOS}_{j,1} = \text{CFCOS}_{j,1} - P_{j,1} \cdot C_i$$

Equation 2

$$\text{ALFA}_K = \text{Re}[\text{ROOT}_K] \quad , \quad \text{BETA}_K = \text{Im}[\text{ROOT}_K]$$

$$\omega_{1,K} = \text{BETA}_K - \text{OMEGA} \quad , \quad \omega_{2,K} = \text{BETA}_K + \text{OMEGA}$$

$$C_{1,K} = \text{ALFA}_K^2 + \omega_{1,K}^2 \quad , \quad C_{2,K} = \text{ALFA}_K^2 + \omega_{2,K}^2$$

$$T_{1,K} = -e^{-\text{ALFA}_K \cdot \text{TO}} [\omega_{1,K} \sin(\omega_{1,K} \cdot \text{TO}) - \text{ALFA}_K \cos(\omega_{1,K} \cdot \text{TO})] / C_{1,K}$$

$$-e^{-\text{ALFA}_K \cdot \text{TO}} [\omega_{2,K} \sin(\omega_{2,K} \cdot \text{TO}) - \text{ALFA}_K \cos(\omega_{2,K} \cdot \text{TO})] / C_{2,K}$$

$$T_{2,K} = -e^{-\text{ALFA}_K \cdot \text{TO}} [-\text{ALFA}_K \sin(\omega_{1,K} \cdot \text{TO}) - \omega_{1,K} \cos(\omega_{1,K} \cdot \text{TO})] / C_{1,K}$$

$$-e^{-\text{ALFA}_K \cdot \text{TO}} [-\text{ALFA}_K \sin(\omega_{2,K} \cdot \text{TO}) - \omega_{2,K} \cos(\omega_{2,K} \cdot \text{TO})] / C_{2,K}$$

$$\text{PR}_{j,1} = \text{AMPL} \cdot \text{Re}[P_{j,1}]$$

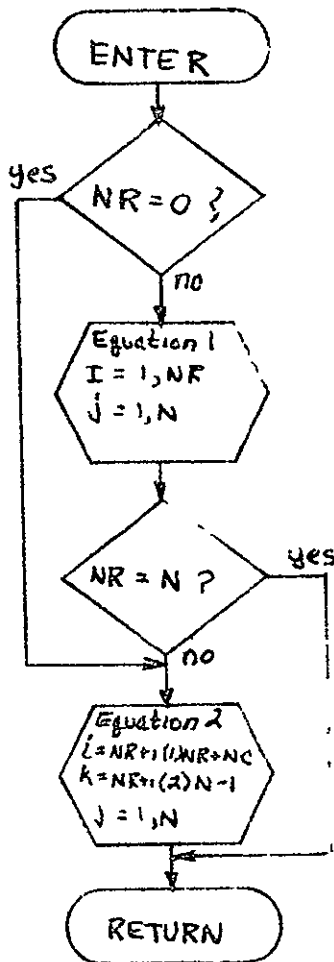
$$\text{PI}_{j,1} = \text{AMPL} \cdot \text{Im}[P_{j,1}]$$

$$\text{STYCOS}_{j,1} = \text{STYCOS}_{j,1} + (-\text{PR}_{j,1} \cdot \omega_{1,K} + \text{PI}_{j,1} \cdot \text{ALFA}_K) / C_{1,K} + (-\text{PI}_{j,1} \cdot \text{ALFA}_K + \text{PR}_{j,1} \cdot \omega_{2,K}) / C_{2,K}$$

$$\text{STYCOS}_{j,2} = \text{STYCOS}_{j,2} + (-\text{PR}_{j,2} \cdot \text{ALFA}_K - \text{PI}_{j,2} \cdot \omega_{1,K}) / C_{1,K} + (-\text{PR}_{j,2} \cdot \text{ALFA}_K - \text{PI}_{j,2} \cdot \omega_{2,K}) / C_{2,K}$$

$$\text{CFCOS}_{j,K} = \text{CFCOS}_{j,K} - T_{1,K} \cdot \text{PI}_{j,1} + T_{2,K} \cdot \text{PR}_{j,1}$$

$$\text{CFCOS}_{j,K+1} = \text{CFCOS}_{j,K+1} + T_{1,K} \cdot \text{PR}_{j,1} + T_{2,K} \cdot \text{PI}_{j,1}$$



SVESIN

Equation 1

$$B_i = \text{ROOT}_i^2 + \text{OMEGA}^2$$

$$C_i = \text{AMPL} \cdot e^{-\text{ROOT}_i \cdot \text{TO}} \left[-\text{ROOT}_i \cdot \sin(\text{OMEGA} \cdot \text{TO}) - \text{OMEGA} \cdot \cos(\text{OMEGA} \cdot \text{TO}) \right] / B_i$$

$$D_i = \text{AMPL} \cdot (-\text{ROOT}_i / B_i) \quad , \quad E_i = \text{AMPL} \cdot (-\text{OMEGA} / B_i)$$

$$\text{STYSIN}_{j,1} = \text{STYSIN}_{j,1} + P_{jL} \cdot D_i$$

$$\text{STYSIN}_{j,2} = \text{STYSIN}_{j,2} + P_{ji} \cdot E_i$$

$$\text{CFSIN}_{j,i} = \text{CFSIN}_{j,i} - P_{ji} \cdot C_i$$

Equation 2

$$\text{ALFA}_k = \text{Re}[\text{ROOT}(k)] \quad , \quad \text{BETA}_k = \text{Im}[\text{ROOT}(k)]$$

$$\omega_{1,k} = \text{BETA}_k - \text{OMEGA} \quad , \quad \omega_{2,k} = \text{BETA}_k + \text{OMEGA}$$

$$C_{1,k} = \text{ALFA}_k^2 + \omega_{1,k}^2 \quad , \quad C_{2,k} = \text{ALFA}_k^2 + \omega_{2,k}^2$$

$$K_{1,k} = -e^{-\text{ALFA}_k \cdot \text{TO}} \left[\omega_{1,k} \cdot \cos(\omega_{1,k} \cdot \text{TO}) + \text{ALFA}_k \sin(\omega_{1,k} \cdot \text{TO}) \right] / C_{1,k} \\ + e^{-\text{ALFA}_k \cdot \text{TO}} \left[\omega_{2,k} \cdot \cos(\omega_{2,k} \cdot \text{TO}) + \text{ALFA}_k \sin(\omega_{2,k} \cdot \text{TO}) \right] / C_{2,k}$$

$$K_{2,k} = -e^{-\text{ALFA}_k \cdot \text{TO}} \left[\omega_{1,k} \cdot \sin(\omega_{1,k} \cdot \text{TO}) - \text{ALFA}_k \cos(\omega_{1,k} \cdot \text{TO}) \right] / C_{1,k} \\ + e^{-\text{ALFA}_k \cdot \text{TO}} \left[\omega_{2,k} \cdot \sin(\omega_{2,k} \cdot \text{TO}) - \text{ALFA}_k \cos(\omega_{2,k} \cdot \text{TO}) \right] / C_{2,k}$$

$$\text{PR}_{ji} = \text{AMPL} \cdot \text{Re}[P_k]$$

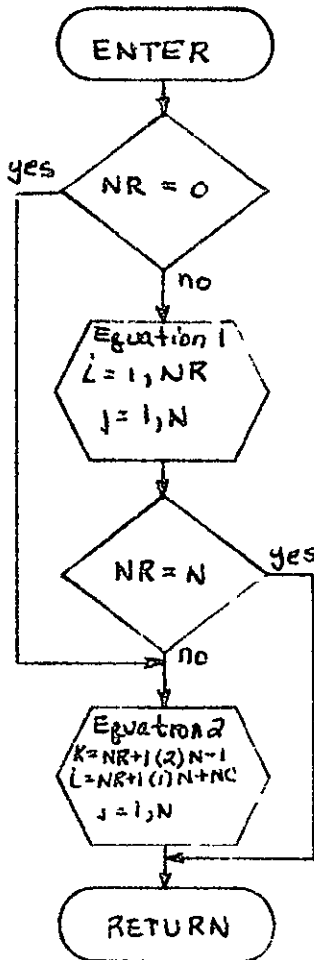
$$\text{PI}_{ji} = \text{AMPL} \cdot \text{Im}[P_k]$$

$$\text{STYSIN}_{j,2} = \text{STYSIN}_{j,2} + (\text{PR}_{ji} \cdot \omega_{1,k} - \text{PI}_{ji} \cdot \text{ALFA}_k) / C_{1,k} + (\text{PI}_{ji} \cdot \text{ALFA}_k - \text{PR}_{ji} \cdot \omega_{2,k}) / C_{2,k}$$

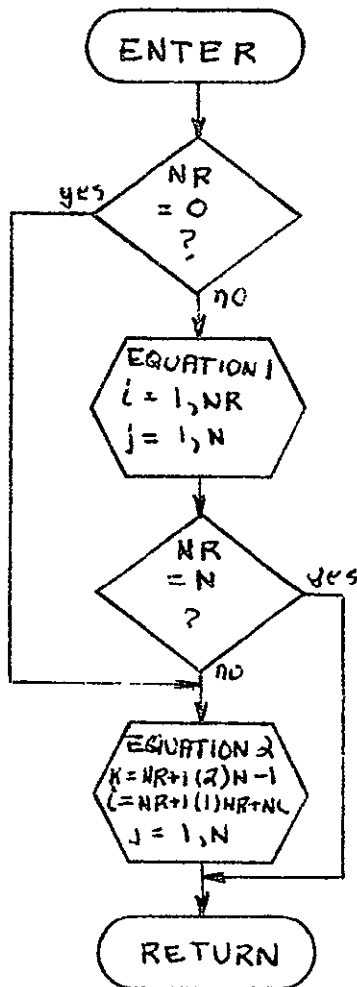
$$\text{STYSIN}_{j,1} = \text{STYSIN}_{j,1} + (-\text{PR}_{ji} \cdot \text{ALFA}_k - \text{PI}_{ji} \cdot \omega_{1,k}) / C_{1,k} + (-\text{PR}_{ji} \cdot \text{ALFA}_k - \text{PI}_{ji} \cdot \omega_{2,k}) / C_{2,k}$$

$$\text{CFSIN}_{j,R+1} = \text{CFSIN}_{j,R+1} + K_{1,k} \cdot \text{PR}_{ji} + K_{2,k} \cdot \text{PI}_{ji}$$

$$\text{CFSIN}_{j,R} = \text{CFSIN}_{j,R} - K_{1,k} \cdot \text{PI}_{ji} + K_{2,k} \cdot \text{PR}_{ji}$$



SVESTP



Equation 1

$$A_L = - \frac{AMPL}{ROOT_L}$$

$$B_L = \frac{AMPL \cdot e^{-ROOT_L \cdot TO}}{ROOT_L}$$

$$COEF_{JL} = COEF_{JL} + P_{JL} \cdot B_L$$

$$STP_J = STP_J + P_{JL} \cdot A_L$$

Equation 2

$$\alpha_K = \text{Re}[ROOT_K]$$

$$\beta_K = \text{Im}[ROOT_K]$$

$$C_K = \frac{2 \cdot AMPL}{\alpha_K^2 + \beta_K^2}$$

$$A_K = C_K \cdot e^{-\alpha_K \cdot TO} [\alpha_K \cos(\beta_K \cdot TO) - \beta_K \sin(\beta_K \cdot TO)]$$

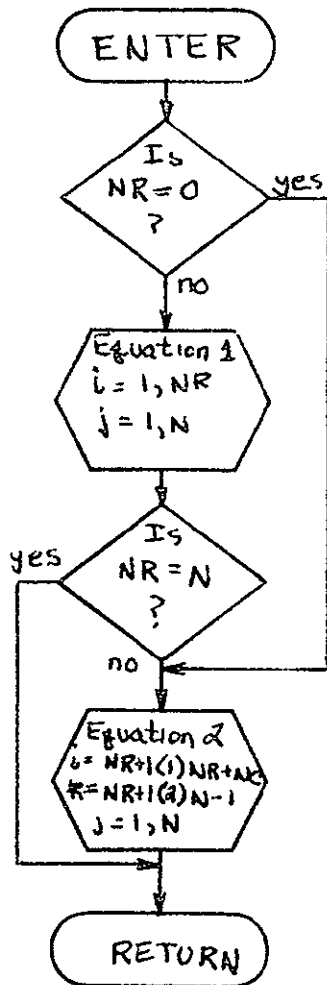
$$B_K = C_K \cdot e^{-\alpha_K \cdot TO} [\beta_K \cos(\beta_K \cdot TO) + \alpha_K \sin(\beta_K \cdot TO)]$$

$$COEF_{J,K} = COEF_{J,K} - \text{Im}[P_{JL}] A_K + \text{Re}[P_{JL}] B_K$$

$$COEF_{J,K+1} = COEF_{J,K+1} + \text{Re}[P_{JL}] A_K + \text{Im}[P_{JL}] B_K$$

$$STP_J = STP_J + C_K [-\alpha_K \text{Re}(P_{JL}) - \beta_K \text{Im}(P_{JL})]$$

INIVAL



Equation 1

$$CF_i = \text{AMPL} \cdot e^{-\text{ROOT}_i \cdot T_0}$$

$$\text{COEFO}_{j,i} = \text{COEFO}_{j,i} + P_{j,i} \cdot CF_i$$

Equation 2

$$\alpha_k = \text{Re} [\text{RCOT}_k]$$

$$\beta_k = \text{Im} [\text{RCOT}_k]$$

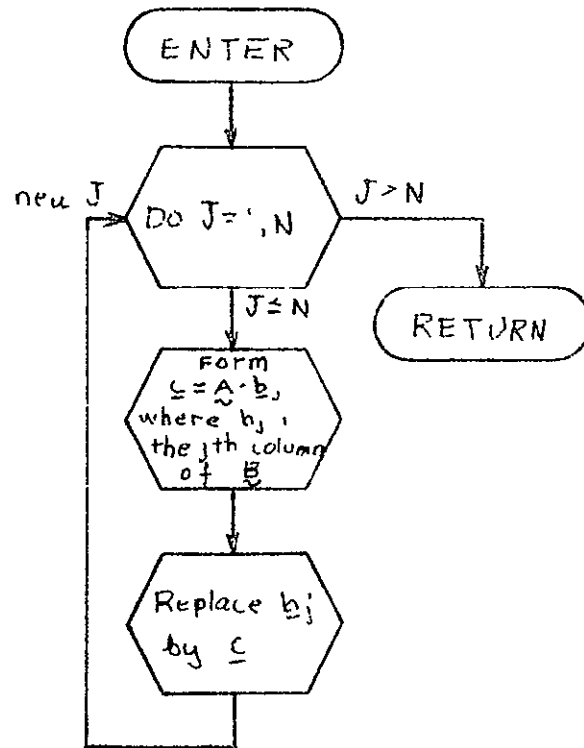
$$A_k = \text{AMPL} \cdot \text{Re} [P_{j,i}]$$

$$B_k = \text{AMPL} \cdot \text{Im} [P_{j,i}]$$

$$\begin{aligned} \text{COEFO}_{j,k} &= \text{COEFO}_{j,k} \\ &+ 2 \cdot e^{-\alpha_k T_0} [A_k \cos(\beta_k T_0) \\ &+ B_k \sin(\beta_k T_0)] \end{aligned}$$

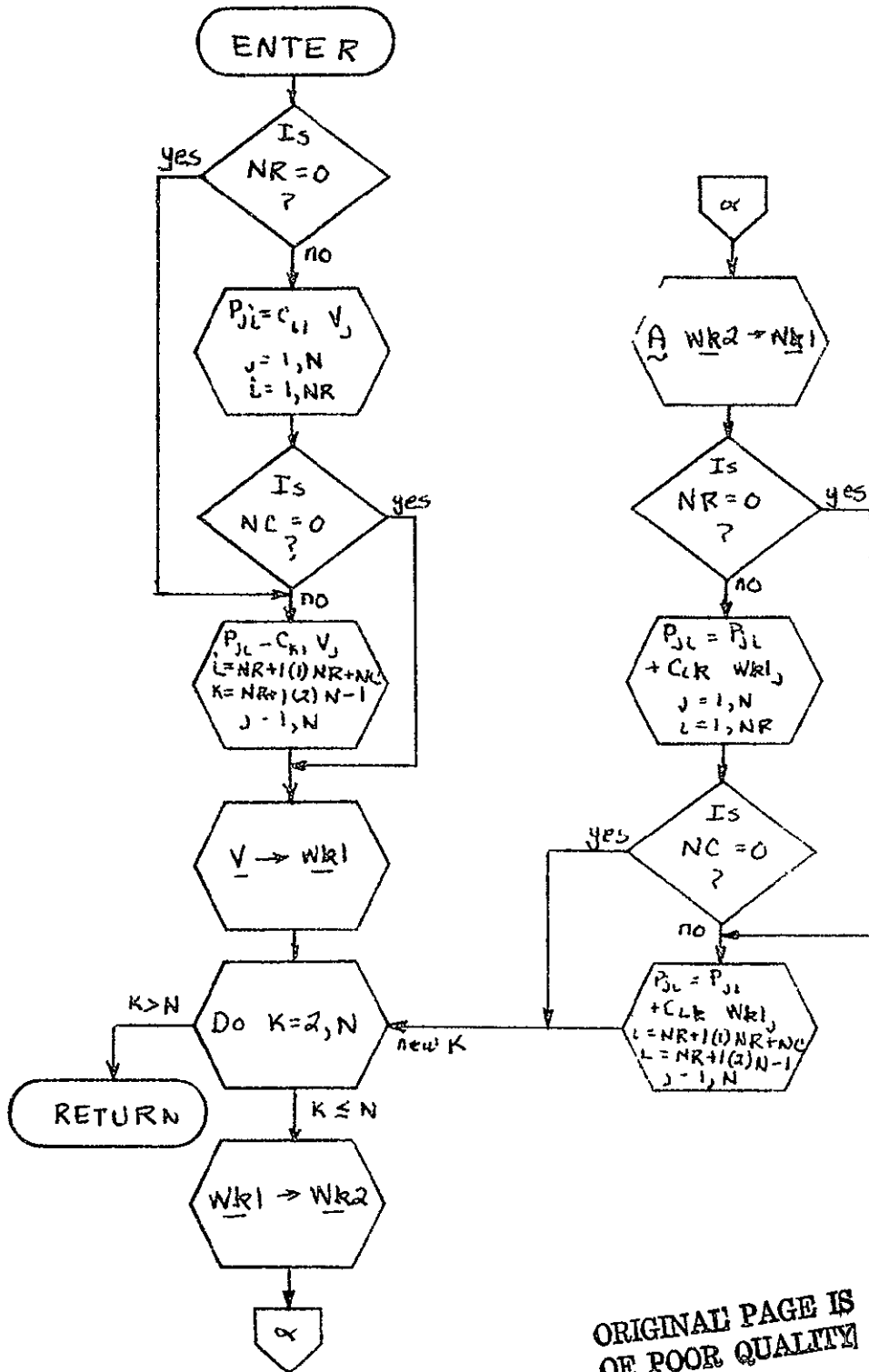
$$\begin{aligned} \text{COEFO}_{j,k+1} &= \text{COEFO}_{j,k+1} \\ &+ 2 \cdot e^{-\alpha_k T_0} [A_k \sin(\beta_k T_0) \\ &+ B_k \cos(\beta_k T_0)] \end{aligned}$$

MPRCD



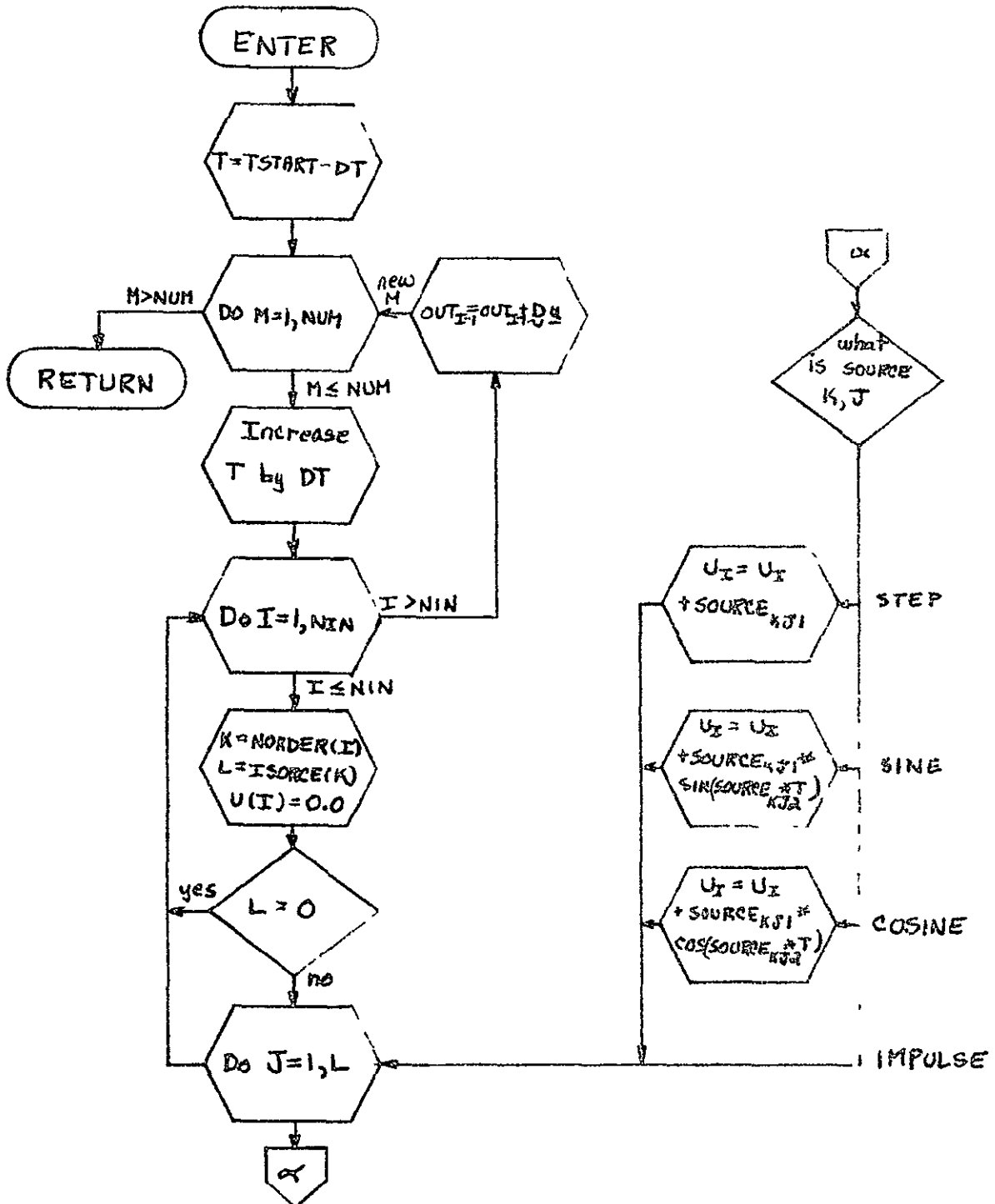
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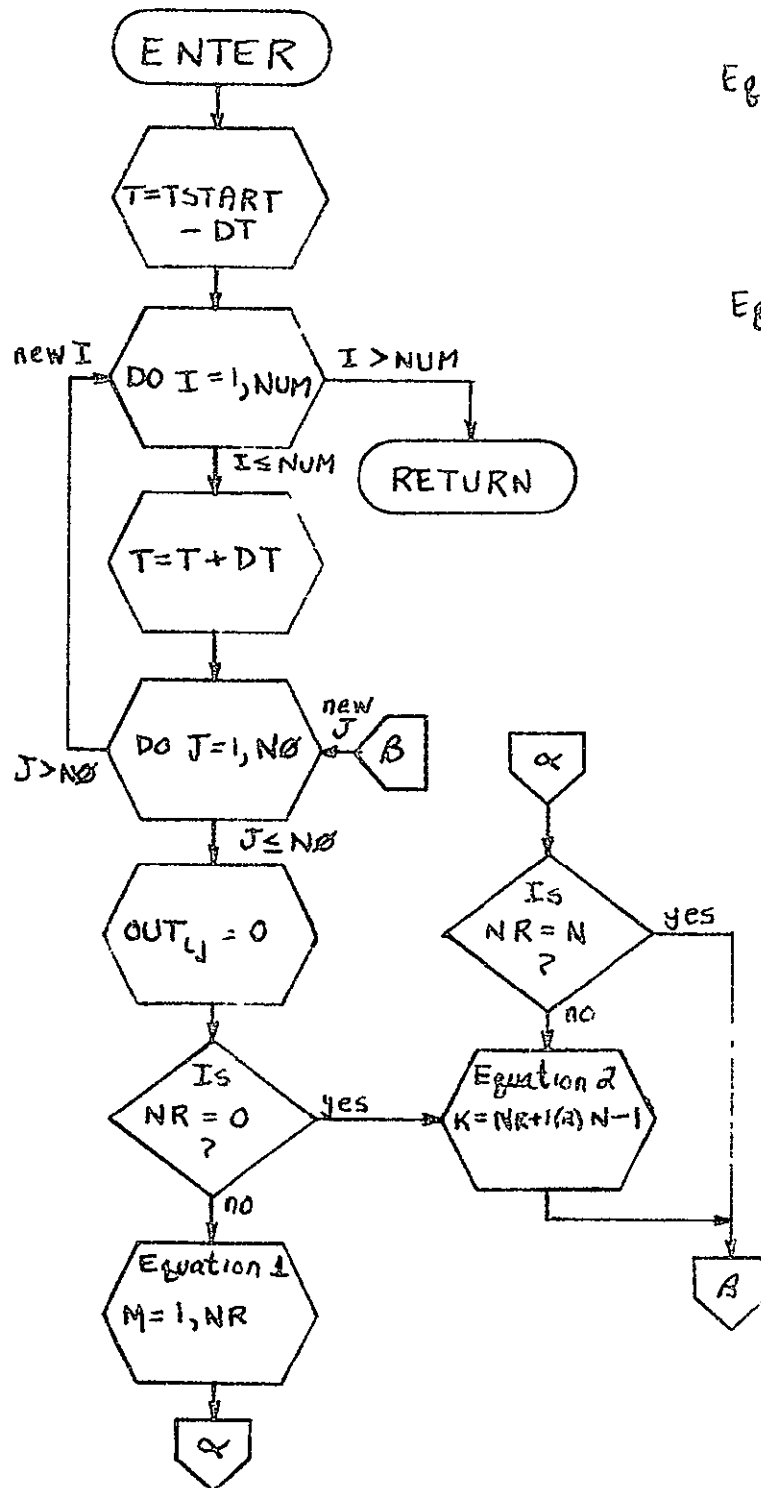


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TRANS



Equation 1

$$OUT_{ij} = OUT_{ij} + COEF_{jm} e^{ROOT_m T}$$

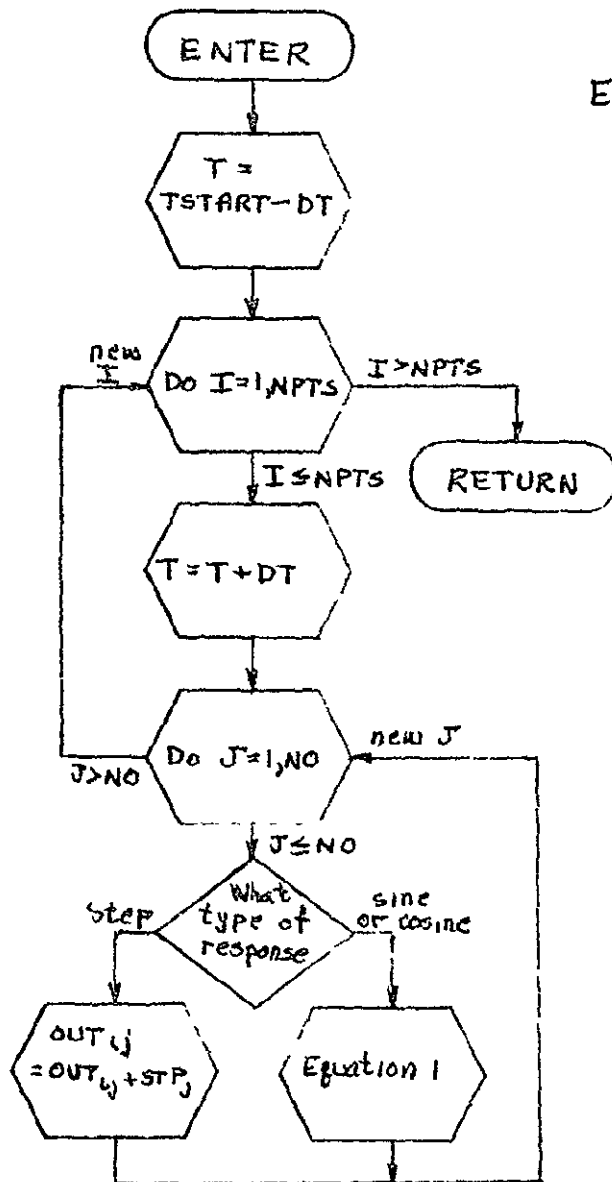
Equation 2

$$\alpha_k = \text{Re} [ROOT_k]$$

$$\beta_k = \text{Im} [ROOT_k]$$

$$OUT_{ij} = CUT_{ij} + [COEF_{jk} \sin(\beta_k T) + COEF_{j,k+1} \cos(\beta_k T)] e^{\alpha_k T}$$

STEADY

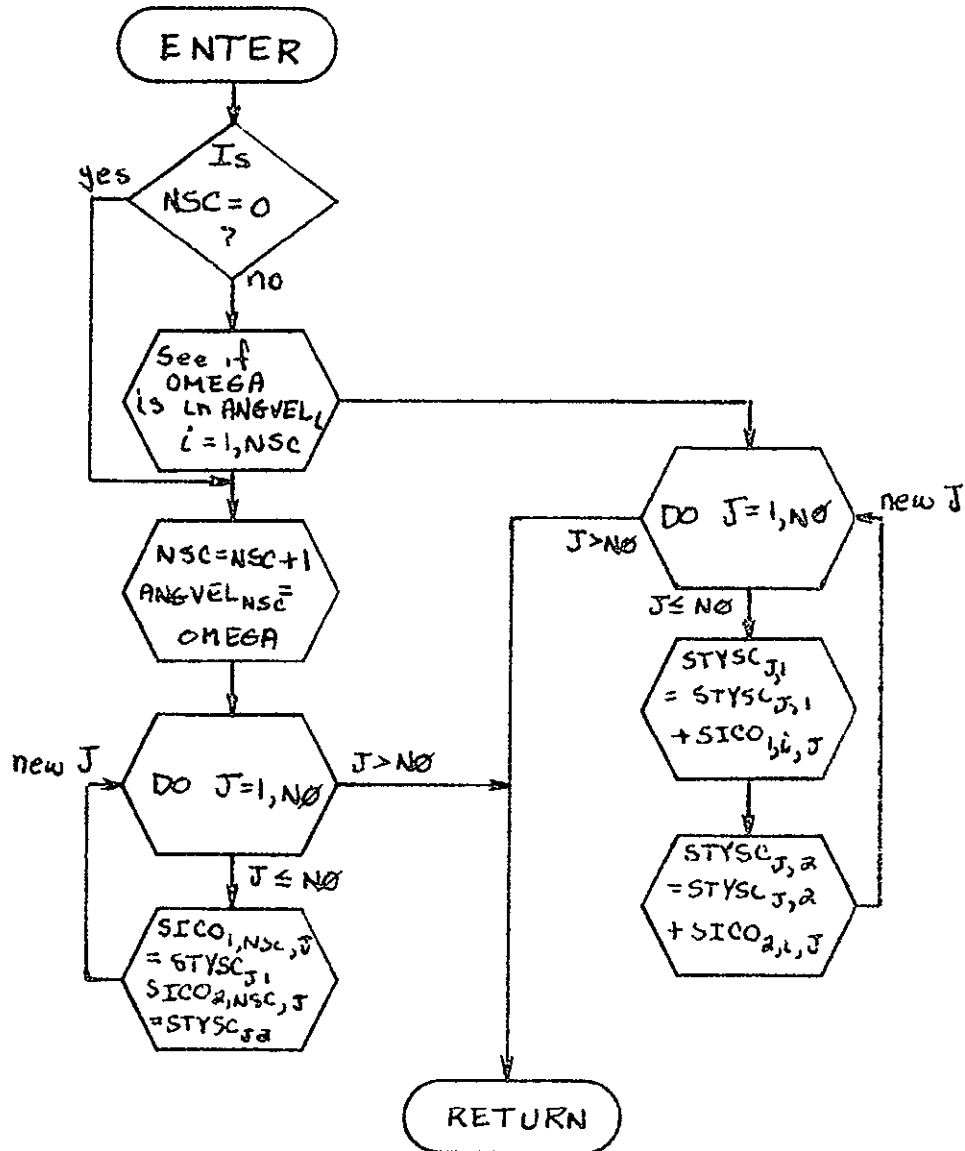


Equation 1

$$OUT_{ij} = OUT_{ij} + STYSC_{j1} \sin(\Omega T) + STYSC_{j2} \cos(\Omega T)$$

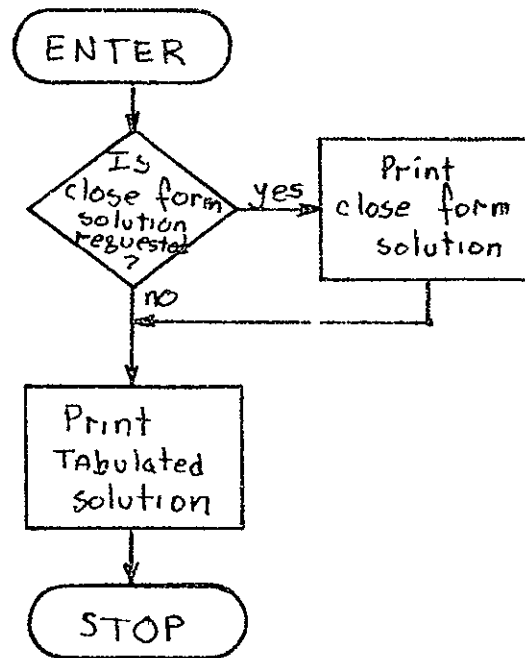
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STORE

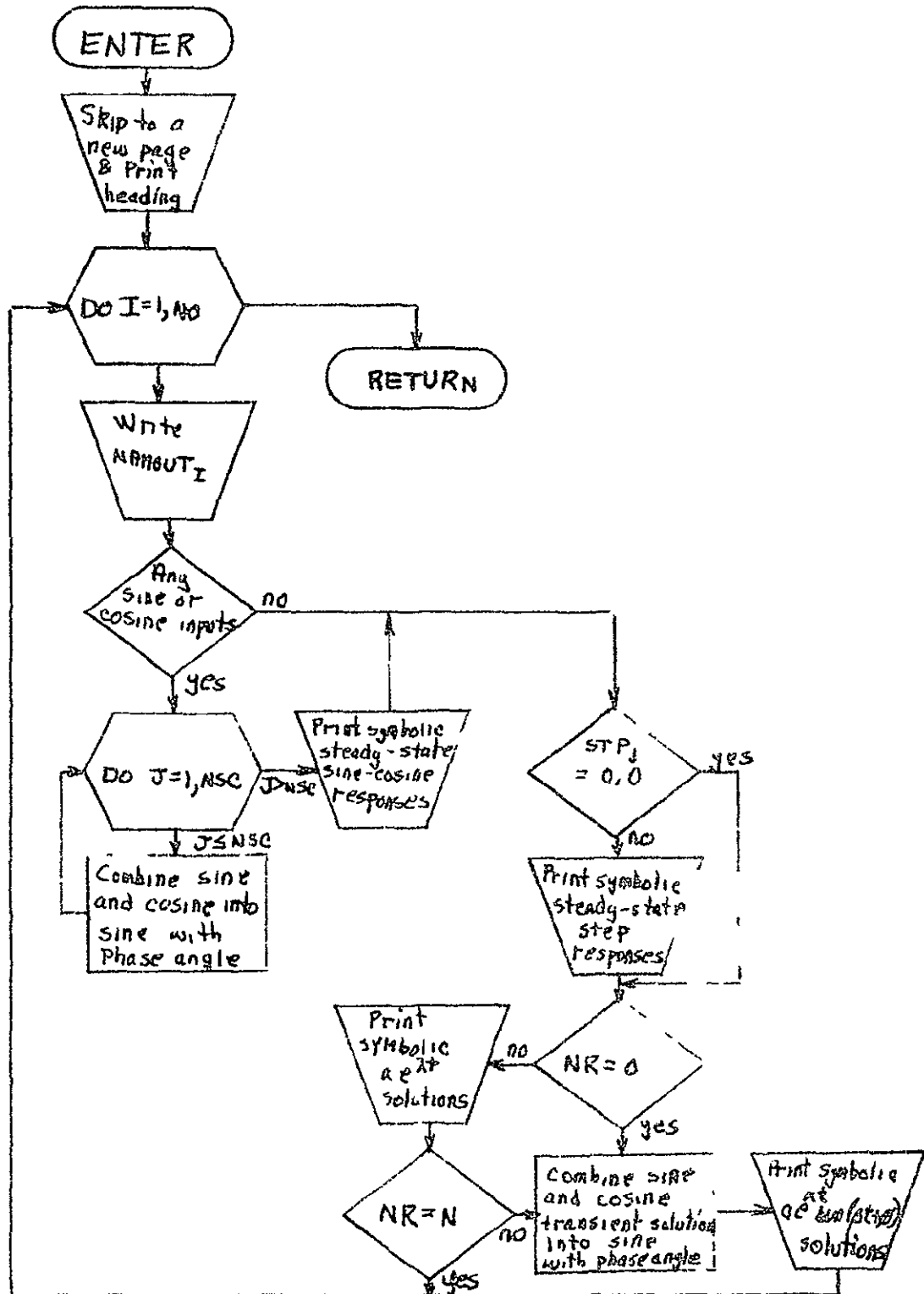


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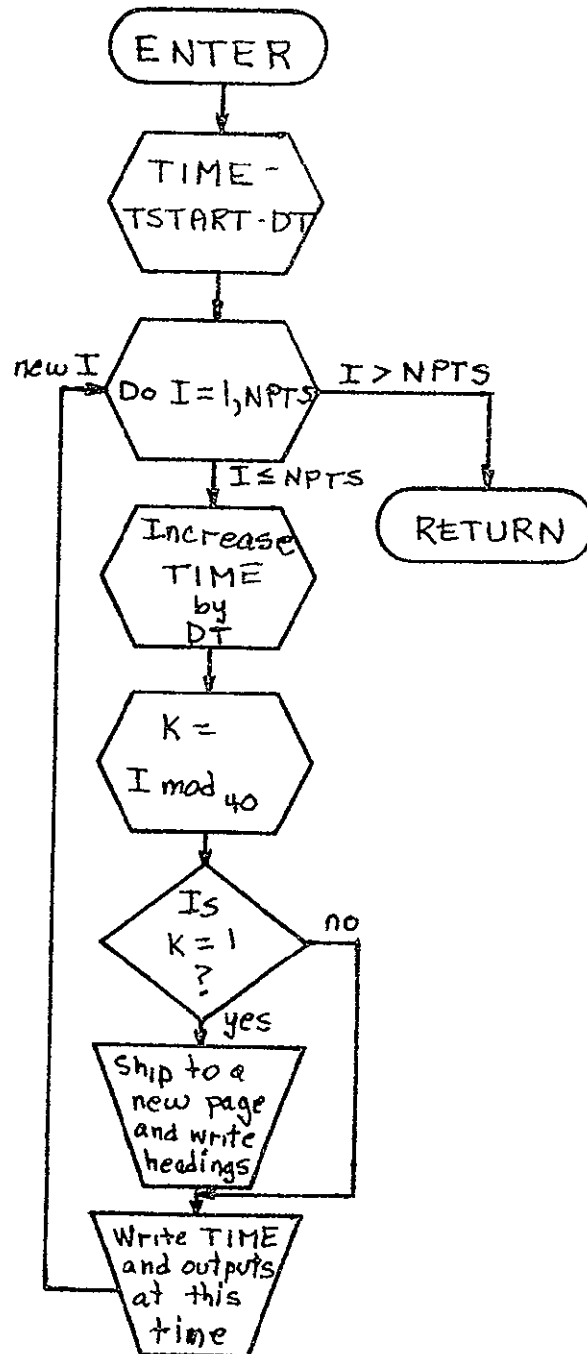
OUT PNT



CLSOL



TABPNT



ATAN3

